



Prepared for:
UTC Aerospace Systems
Rockford, IL

Prepared by:
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December 2015

Third Quarter 2015 GMZ Monitoring and System Performance Report

UTC Aerospace Systems Plants 1/2 Facility
Area 9/10 Remedial Action
Southeast Rockford Groundwater Contamination
Superfund Site
2421 11th Street
Rockford, IL 61104
ILD 981000417

US EPA RECORDS CENTER REGION 5



521806



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Letter of Transmittal

Attention:	Mr. Timothy Drexler (USEPA) Mr. Brian Conrath (IEPA)	Date:	December 11, 2015
<p>UTC Aerospace Systems Plants 1/2 Facility Area 9/10 Remedial Action Southeast Rockford Groundwater Contamination Superfund Site 2421 11th Street Rockford, Illinois 61104</p>			
Project reference:	ILD981000417	Project number:	60339110

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Enclosed please find the Third Quarter 2015 GMZ Monitoring and System Performance Report for UTC Aerospace Systems Plants 1/2 Facility, Area 9/10 Remedial Action, Southeast Rockford Groundwater Contamination Superfund Site, Rockford, Illinois.

Thank You.

Peter Hollatz, P.E.

cc: Mr. Scott Moyer, United Technologies Corporation (cd only)
 Ms. Diane Bellantoni, UTC Aerospace Systems (cd only)
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Subject: Third Quarter 2015 GMZ Monitoring and System Performance Report
UTC Aerospace Systems Plants 1/2 Facility
Area 9/10 Remedial Action
Southeast Rockford Groundwater Contamination Superfund Site
2421 11th Street
Rockford, Illinois 61104
ILD981000417
AECOM Project No. 60339110

Dear Messrs. Drexler and Conrath:

This Quarterly Groundwater Management Zone (GMZ) Monitoring and System Performance Report has been prepared by AECOM Technical Services Inc. (AECOM) on behalf of UTC Aerospace Systems (UTAS, fka Hamilton Sundstrand or HS). In accordance with the approved March 2007 Operation, Maintenance, and Monitoring Plan (OM&M Plan) and the United States Environmental Protection Agency (EPA) letter dated April 15, 2011 providing approval for combining project reporting documents, this report contains a summary of the following: 1) GMZ groundwater monitoring data; 2) the Phase 1 and Phase 2 air sparge/soil vapor extraction (AS/SVE) system performance data; 3) the Phase 1 and Phase 2 AS/SVE system process air analytical data; 4) GMZ wells that contain contaminants of concern (COCs) above Preliminary Remediation Goals (PRGs); and 5) Quarterly Progress Report for Fourth Quarter 2015.

As approved in the April 15, 2011 letter from Timothy Drexler, interpretation of collected groundwater quality and system performance data will be included in the Annual GMZ Monitoring and System Performance Report submitted in March of the subsequent year. This quarterly report provides the current environmental data including: tables and figures summarizing the results of third quarter 2015 GMZ monitoring and AS/SVE system performance data, supporting field data sheets and laboratory analytical reports, and the Quarterly Progress Report covering the period from September 1, 2015, to November 30, 2015.

The objective of AS/SVE system operation is to treat leachate-impacted groundwater at the UTAS Plants 1/2 (Site) property. The implemented remedy was specifically targeted to address an area of the Site where COCs were originally present in leachate/groundwater at concentrations that were two or more orders of magnitude greater than their PRGs. Though the treatment area was not fully defined when the 2002 Record of Decision (ROD) for Operable Unit 3 (OU3) was issued, the entire Site was identified/defined in the ROD as a "source location" within the larger established "Source Area 9/10" (Area 9/10) based on data collected prior to the ROD¹. The ROD further required that the Site remedy include the establishment of a GMZ for this "source location" (the Site) whose lateral limits were consequently defined as the Site property boundaries. Two Site GMZs, GMZ 1 (Site property north of railroad tracks) and GMZ 2 (Site property south of railroad tracks), were approved by the Illinois EPA in 2008. Monitoring wells within the Site GMZs are routinely sampled, and the groundwater analytical results are compared to OU3 PRGs to evaluate the effectiveness of the remedy.

During the third quarter 2015 reporting period, the following six GMZ well locations along the Site boundary contained COCs at concentrations above PRGs:

GMZ Monitoring Well ID	COC ^[1] Concentrations > PRG (Increase (+) or Decrease (-) from Previous Quarter)
GMZ01	PCE (+)
SMW04	cis-1,2-DCE (-); PCE (+); VC (-)
SMW08	cis-1,2-DCE (+); PCE (+)
SMW19	TCE (-)
PMW01	PCE (+)
PMW02	PCE (-)

^[1] Trichloroethene (TCE), cis-1,2-Dichloroethene (cis-1,2-DCE), Tetrachloroethene (PCE), Vinyl Chloride (VC)

The above-noted decreases/increases in concentrations represent a relative change in COC concentrations (above the PRG) between the two most recent quarters of data. Such changes should not be viewed as an indication of a trend without further statistical evaluation.

While PRGs are used to assess on-going remedy effectiveness at the Site, the continued operation of the AS/SVE remedy will be dependent on the attainment of Alternate Cleanup Levels (ACLs) at the downgradient Site GMZ boundary. COC ACLs have not yet been established/approved for the Site, but the ACLs will represent the maximum allowable concentration at the Site boundary that will not result in a COC exceedance of a PRG at the Area 9/10 boundary downgradient of the Site.

¹ See EPA Superfund Record of Decision Southeast Rockford Ground Water Contamination, 2002. EPA/ROD/R05-02/077 2002.

Achieving ACLs at the downgradient Site boundary will demonstrate that the Site is protective of human or environmental receptors at the downgradient Area 9/10 boundary, and that continued active remediation is no longer warranted. The downgradient Area 9/10 boundary is located at Harrison Avenue to the south and 6th Street to the west.

The formulation of ACLs is consistent with the attainment of the OU3 ROD Remedial Action Objective (RAO) for groundwater specified in the ROD² and the objectives analysis/Remedial Action Process Flow Diagram (RAPFD) developed and approved for use by the EPA and Illinois EPA at the Site. The RAPFD, and the conditions for the performance of an objectives analysis and use of ACLs at the Site, are provided in the Statement of Work attached to the UTAS (fka Hamilton Sundstrand) facility Consent Decree³ and included in subsequent approved Remedial Design documents for the Site.

Please contact either of the undersigned with any questions you may have on the information provided.

Prepared by:



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² The OU3 ROD RAO for groundwater media is to: "Prevent the further migration of contamination from the source area that would result in degradation of site-wide groundwater or surface water to levels in excess of state or federal standards, or that pose a threat to human health or the environment."

³ See the Statement of Work in Appendix C of the Consent Decree between Hamilton Sundstrand Corporation and the United States Environmental Protection Agency (Civil Action Number 08 C 50129), Section II.D.2, *Implementation of Remedial Action and Attainment of Performance Standards* (pages 9 and 10).

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Tables

Table 1
Fourth Quarter 2014 to Third Quarter 2015 Cumulative Groundwater Elevations
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

Well ID	Top of Casing Elevation (ft)	Depth to Groundwater (ft BTOC)	Groundwater Elevaion (ft AMSL)	Depth to Groundwater (ft BTOC)	Groundwater Elevation (ft AMSL)	Depth to Groundwater (ft BTOC)	Groundwater Elevation (ft AMSL)	Depth to Groundwater (ft BTOC)	Groundwater Elevaion (ft AMSL)
		12/8/2014		2/16/2015		5/11/2015		8/4/2015	
MW07FGA	727.49	28.78	698.71	29.17	698.32	29.47	698.02	29.33	698.16
MW203	728.58	29.62	698.96	29.96	698.62	30.30	698.28	30.12	698.46
SMW01	729.71	31.64	698.07	32.00	697.71	32.26	697.45	32.08	697.63
SMW02	726.77	28.11	698.66	28.54	698.23	28.77	698.00	28.59	698.18
SMW04	728.51	30.75	697.76	31.15	697.36	31.41	697.10	31.22	697.29
SMW08	728.81	30.97	697.84	31.34	697.47	31.64	697.17	31.42	697.39
SMW19	728.49	29.79	698.70	30.12	698.37	30.44	698.05	30.31	698.18
SMW20	727.69	29.94	697.75	30.26	697.43	30.52	697.17	30.41	697.28
SMW21	727.25	29.30	697.95	29.71	697.54	30.02	697.23	29.88	697.37
GMZ01	731.41	33.56	697.85	33.95	697.46	34.19	697.22	34.02	697.39
GMZ02	728.76	31.06	697.70	31.47	697.29	31.69	697.07	31.56	697.20
GMZ03	728.22	30.47	697.75	29.91	698.31	31.08	697.14	30.95	697.27
GMZ04	726.84	28.73	698.11	29.18	697.66	29.41	697.43	29.27	697.57
BGW01	728.19	29.65	698.54	30.01	698.18	30.29	697.90	30.08	698.11
BGW02	728.81	30.01	698.80	30.46	698.35	30.77	698.04	30.57	698.24
BGW03	728.96	30.12	698.84	30.48	698.48	30.79	698.17	30.66	698.30
RAMW01	728.91	31.13	697.78	31.54	697.37	31.75	697.16	31.60	697.31
RAMW02	728.90	31.00	697.90	31.40	697.50	31.64	697.26	31.48	697.42
RAMW03	728.71	30.80	697.91	31.18	697.53	31.45	697.26	31.29	697.42
RAMW04	728.80	30.63	698.17	31.04	697.76	31.32	697.48	31.14	697.66
RAMW05	727.65	29.50	698.15	29.94	697.71	30.20	697.45	30.02	697.63
RAMW06	727.64	29.53	698.11	30.01	697.63	30.22	697.42	30.07	697.57
RAMW07	732.20	34.00	698.20	34.41	697.79	34.68	697.52	34.53	697.67
RAMW08	728.45	30.14	698.31	30.54	697.91	30.83	697.62	30.64	697.81
PMW01	728.88	31.26	697.62	31.64	697.24	31.79	697.09	31.71	697.17
PMW02	728.88	31.20	697.68	31.58	697.30	31.78	697.10	31.64	697.24
Ave. GW Elev. (ft AMSL)		698.15		697.79		697.49		697.65	

Notes:

NM = Not monitored

ft = feet

ft BTOC = feet below top of casing

ft AMSL = feet above mean sea level

All site well top of casing elevations re-surveyed on May 24, 2011.

RAMW04 riser was lowered due to ice damage that occurred during the 2013 winter. Well was resurveyed on July 1, 2013.

Table 2
Fourth Quarter 2014 to Third Quarter 2015 Groundwater Analytical Results - GMZ Wells
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

				Trichloroethene (TCE)	Methylene Chloride (Dichloromethane)	1,1-Dichloroethene	1,1-Dichloroethane	1,2-Dichloroethane	cis-1,2-Dichloroethene	trans-1,2-Dichloroethene	1,1,1-Trichloroethane	1,1,2-Trichloroethane	Ethylbenzene	Tetrachloroethene (PCE)	Toluene	Vinyl chloride	
Preliminary Remediation Goals (PRG) ^A				0.005 ^A _c	0.005 ^A _c	0.007 ^{b,c} ^A	0.7 ^A	0.005 ^A _c	0.07 ^A _c	0.1 ^A _c	0.2 ^{b,c} ^A	0.005 ^A _c	0.7 ^A _c	0.005 ^A _c	1.0 ^A _c	0.002 ^A _c	
Well	Sample ID	Sample Date	Sample Type	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	
GMZ01	HS SER-GMZ01-120814	8-Dec-14		0.0014	0.0020 U	0.0010 U	0.0075	0.0010 U	0.0044	0.0010 U	0.0034	0.0010 U	0.0010 U	0.0126 ^A	0.0010 U	0.0010 U	
	HS SER-GMZ01-021715	17-Feb-15		0.0012	0.0020 U	0.0010 U	0.0065	0.0010 U	0.0020	0.0010 U	0.0042	0.0010 U	0.0010 U	0.0044	0.0010 U	0.0010 U	
	HS SER-GMZ01-051115	11-May-15		0.0019	0.0020 U	0.0010 U	0.0088	0.0010 U	0.0022	0.0010 U	0.0063	0.0010 U	0.0010 U	0.0157 ^A	0.0010 U	0.0010 U	
	HS SER-GMZ01-080415	4-Aug-15		0.0018	0.0020 U	0.0010 U	0.0066	0.0010 U	0.0037	0.0010 U	0.0064	0.0010 U	0.0010 U	0.0183 ^A	0.0010 U	0.0010 U	
GMZ02	HS SER-GMZ02-120914	9-Dec-14		0.00037 J	0.0020 U	0.0010 U	0.0016	0.0010 U	0.00073 J	0.0010 U	0.0065	0.0010 U	0.0010 U	0.00064 J	0.0010 U	0.0010 U	
	HS SER-GMZ02-021915	19-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.0028	0.0010 U	0.00074 J	0.0010 U	0.0039	0.0010 U	0.0010 U	0.00038 J	0.0010 U	0.0010 U	
	HS SER-GMZ02-051315	13-May-15		0.0010 U	0.0020 U	0.0010 U	0.0018	0.0010 U	0.00042 J	0.0010 U	0.0041	0.0010 U	0.0010 U	0.00057 J	0.0010 U	0.0010 U	
	HS SER-GMZ-080515	5-Aug-15		0.00040 J	0.0020 U	0.0010 U	0.0016	0.0010 U	0.00057 J	0.0010 U	0.0030	0.0010 U	0.0010 U	0.00043 J	0.0010 U	0.0010 U	
GMZ03	HS SER-GMZ03-120914	9-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0098 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-DUP01-120914	9-Dec-14	Field Duplicate	0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0093 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-GMZ03-021815	18-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.00037 J	0.0010 U	0.0010 U	0.0010 U	0.0010 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-GMZ03-021815	18-Feb-15	Field Duplicate	0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00090 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-GMZ03-051315	13-May-15		0.0010 U	0.0020 U	0.0010 U	0.00076 J	0.0010 U	0.00050 J	0.0010 U	0.00090 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-DUP01-051315	13-May-15	Field Duplicate	0.0010 U	0.0020 U	0.0010 U	0.00078 J	0.0010 U	0.00055 J	0.0010 U	0.00089 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-GMZ03-080515	5-Aug-15		0.0010 U	0.0020 U	0.0010 U	0.00028 J	0.0010 U	0.0010 U	0.00035 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-DUP01-080515	5-Aug-15	Field Duplicate	0.0010 U	0.0020 U	0.0010 U	0.00026 J	0.0010 U	0.0010 U	0.00036 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
GMZ04	HS SER-GMZ04-121014	10-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00065 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-GMZ04-021815	18-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.00038 J	0.0010 U	0.0023	0.0010 U	0.0036	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-GMZ04-051215	12-May-15		0.0010 U	0.0020 U	0.0010 U	0.0024	0.0010 U	0.0030	0.0010 U	0.0113	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-GMZ04-080515	5-Aug-15		0.0026	0.0020 U	0.00093 J	0.0011	0.0010 U	0.0035	0.0010 U	0.0536	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
MW07FGA	HS SER-MW07FGA-120914	9-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00043 J	0.0010 U	0.0010 U	0.0010 U	0.0012	0.0010 U	0.0010 U	
	HS SER-MW07FGA-021715	17-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00044 J	0.0010 U	0.0010 U	0.0010 U	0.00076 J	0.0010 U	0.0010 U	
	HS SER-MW07FGA-051215	12-May-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00046 J	0.0010 U	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U	
	HS SER-MW07FGA-080515	5-Aug-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U	
MW203	HS SER-MW203-120914	9-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0044	0.0010 U	0.0010 U	
	HS SER-MW203-021815	18-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0036	0.0010 U	0.0010 U	
	HS SER-MW203-051115	11-May-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0038	0.0010 U	0.0010 U	
	HS SER-MW203-080415	4-Aug-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0037	0.0010 U	0.0010 U	
SMW01	HS SER-SMW01-120914	9-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0012	0.0010 U	0.0010 U	0.0012	0.0010 U	0.0010 U	
	HS SER-SMW01-021715	17-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00078 J	0.0010 U	0.0010 U	0.0010 U	0.00097 J	0.0010 U	0.0010 U	
	HS SER-SMW01-051215																

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UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

				Trichloroethene (TCE)	Methylene Chloride (Dichloromethane)	1,1-Dichloroethene	1,1-Dichloroethane	1,2-Dichloroethane	cis-1,2-Dichloroethene	trans-1,2-Dichloroethene	1,1,1-Trichloroethane	1,1,2-Trichloroethane	Ethylbenzene	Tetrachloroethene (PCE)	Toluene	Vinyl chloride
Preliminary Remediation Goals (PRG) ^A				0.005 ^a	0.005 ^a	0.007 ^{b,c}	0.7 ^a	0.005 ^a	0.07 ^a	0.1 ^a	0.2 ^{b,c}	0.005 ^a	0.7 ^a	0.005 ^a	1.0 ^a	0.002 ^a
Well	Sample ID	Sample Date	Sample Type	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L
SMW08	HS SER-SMW08-120814	8-Dec-14		0.0039	0.0020 U	0.0010 U	0.0165	0.0010 U	0.123 ^a	0.00088 J	0.0067	0.0010 U	0.0010 U	0.0227 ^a	0.0010 U	0.0010 U
	HS SER-SMW08-021715	17-Feb-15		0.0042	0.0020 U	0.0010 U	0.0103	0.0010 U	0.125 ^a	0.00085 J	0.0016	0.0010 U	0.0010 U	0.0092 ^a	0.0010 U	0.0010 U
	HS SER-SMW08-051115	11-May-15		0.0021	0.0020 U	0.0010 U	0.0084	0.0010 U	0.0040	0.0010 U	0.0097	0.0010 U	0.0010 U	0.0199 ^a	0.0010 U	0.0010 U
	HS SER-SMW08-080415	4-Aug-15		0.0021	0.0020 U	0.0011	0.0165	0.0010 U	0.140 ^a	0.0013	0.0040	0.0010 U	0.0010 U	0.0243 ^a	0.0010 U	0.0010 U
SMW19	HS SER-SMW19-121014	10-Dec-14		0.0194 ^a	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U
	HS SER-SMW19-021915	19-Feb-15		0.0177 ^a	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0019	0.0010 U	0.00041 J	0.0010 U	0.0010 U	0.00091 J	0.0010 U	0.0010 U
	HS SER-SMW19-051215	12-May-15		0.0145 ^a	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0011	0.0010 U	0.00034 J	0.0010 U	0.0010 U	0.0010	0.0010 U	0.0010 U
	HS SER-SMW19-080515	5-Aug-15		0.0106 ^a	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0007 J	0.0010 U	0.00034 J	0.0010 U	0.0010 U	0.00086 J	0.0010 U	0.0010 U
SMW20	HS SER-SMW20-121014	10-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-SMW20-021815	18-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-SMW20-051215	12-May-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.00043 J	0.0010 U	0.00027 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-SMW20-080515	5-Aug-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
SMW21	HS SER-SMW21-121014	10-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.00036 J	0.0010 U	0.0031	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-SMW21-021815	18-Feb-15		0.00037 J	0.0020 U	0.0010 U	0.00047 J	0.0010 U	0.0028	0.0010 U	0.0156	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-SMW21-051215	12-May-15		0.00022 J	0.0020 U	0.00069 J	0.00098 J	0.0010 U	0.0054	0.0010 U	0.0315	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-SMW21-080515	5-Aug-15		0.00084 J	0.0020 U	0.0010 U	0.00025 J	0.0010 U	0.00099 J	0.0010 U	0.0100	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
PMW01	HS SER-PMW01-120914	9-Dec-14		0.0012	0.0020 U	0.00054 J	0.0057	0.0010 U	0.00063 J	0.0010 U	0.0259	0.0010 U	0.0010 U	0.0177 ^a	0.0010 U	0.0010 U
	HS SER-PMW01-021815	18-Feb-15		0.0011	0.0020 U	0.0010 U	0.0052	0.0010 U	0.00070 J	0.0010 U	0.0202	0.0010 U	0.0010 U	0.0185 ^a	0.0010 U	0.0010 U
	HS SER-PMW01-051315	13-May-15		0.0014	0.0020 U	0.0010	0.0088	0.0010 U	0.0016	0.0010 U	0.0394	0.0010 U	0.0010 U	0.0178 ^a	0.0010 U	0.0010 U
	HS SER-PMW01-080615	6-Aug-15		0.0013	0.0020 U	0.0018	0.0086	0.0010 U	0.0011	0.0010 U	0.0547	0.0010 U	0.0010 U	0.0192 ^a	0.0010 U	0.0010 U
PMW02	HS SER-PMW02-120914	9-Dec-14		0.0028	0.0020 U	0.00060 J	0.0074	0.0010 U	0.0096	0.0010 U	0.0056	0.0010 U	0.0010 U	0.0144 ^a	0.0010 U	0.0012
	HS SER-PMW02-021815	18-Feb-15		0.0020	0.0020 U	0.0010 U	0.0036	0.0010 U	0.0011	0.0010 U	0.0162	0.0010 U	0.0010 U	0.0210 ^a	0.0010 U	0.0010 U
	HS SER-PMW02-051315	13-May-15		0.0010	0.0020 U	0.0010 U	0.0054	0.0010 U	0.0055	0.0010 U	0.0036	0.0010 U	0.0010 U	0.0149 ^a	0.0010 U	0.0021 ^a
	HS SER-PMW02-080615	6-Aug-15		0.0018	0.0020 U	0.0010 U	0.0069	0.0010 U	0.0043	0.0010 U	0.0042	0.0010 U	0.0010 U	0.0088 ^a	0.0010 U	0.0016

Notes:

PRG Preliminary Remediation Goals (PRGs) from the Record of Decision (ROD)

^A Class 1 - Groundwater Remediation Objectives

6.5^a Concentration exceeds the indicated standard.

15.2 Concentration was detected but did not exceed applicable standards.

0.03 U The analyte was not detected above the laboratory estimated quantitation limit.

0.50 U Laboratory estimated quantitation limit exceeded standard.

n/v No standard/guideline value.

- Parameter not analyzed / not available.

mg/L milligrams per liter

^{b,c} Oral Reference Dose and/or Reference Concentration under review by USEPA. Listed values subject to change.

Value listed is also the Groundwater Quality Standard for this chemical pursuant to 35 Ill.Adm.Code 620.410 for

Class I Groundwater or 35 Ill.Adm.Code 620.420 for Class II Groundwater.

^c Value listed is also the Groundwater Quality Standard for this chemical pursuant to 35 Ill.Adm.Code 620.410 for

Class I Groundwater or 35 Ill.Adm.Code 620.420 for Class II Groundwater.

* LCS or LCSD exceeds the control limits

B The analyte was detected in the method, field and/or trip blank.

H Sample was prepped or analyzed beyond the specified holding time

J Indicates estimated value.

NJ The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

Table 3
Fourth Quarter 2014 to Third Quarter 2015 Groundwater Analytical Results - Performance Wells
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

				Trichloroethene (TCE)	Methylene Chloride (Dichloromethane)	1,1-Dichloroethene	1,1-Dichloroethane	1,2-Dichloroethane	cis-1,2-Dichloroethene	trans-1,2-Dichloroethene	1,1,1-Trichloroethane	1,1,2-Trichloroethane	Ethylbenzene	Tetrachloroethene (PCE)	Toluene	Vinyl chloride
Preliminary Remediation Goals (PRG) ^A				0.005 ^a	0.005 ^a	0.007 ^{b,c} ^a	0.7 ^a	0.005 ^c ^a	0.07 ^c ^a	0.1 ^c ^a	0.2 ^{b,c} ^a	0.005 ^c ^a	0.7 ^c ^a	0.005 ^c ^a	1.0 ^c ^a	0.002 ^c ^a
Well	Sample ID	Sample Date	Sample Type	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L
RAMW01	HS SER-RAMW01-121014	10-Dec-14		0.0011	0.0020 U	0.0010 U	0.0043	0.0010 U	0.00057 J	0.0010 U	0.0042	0.0010 U	0.0010 U	0.0073 ^a	0.0010 U	0.0010 U
	HS SER-RAMW01-021715	17-Feb-15		0.00095 J	0.0020 U	0.0010 U	0.0040	0.0010 U	0.00055 J	0.0010 U	0.0029	0.0010 U	0.0010 U	0.0049	0.0010 U	0.0010 U
	HS SER-RAMW01-051415	14-May-15		0.0010	0.0020 U	0.0010 U	0.0030	0.0010 U	0.0010	0.0010 U	0.0029	0.0010 U	0.0010 U	0.0047	0.0010 U	0.0010 U
	HS SER-RAMW01-080615	6-Aug-15		0.0012	0.0020 U	0.0010 U	0.0037	0.0010 U	0.00070 J	0.0010 U	0.0039	0.00022 J	0.0010 U	0.0052 ^a	0.0010 U	0.0010 U
RAMW02	HS SER-RAMW02-121014	10-Dec-14		0.00063 J	0.0020 U	0.0010 U	0.0033	0.0010 U	0.00033 J	0.0010 U	0.0044	0.0010 U	0.0010 U	0.0035	0.0010 U	0.0010 U
	HS SER-RAMW02-021715	17-Feb-15		0.00055 J	0.0020 U	0.0010 U	0.0037	0.0010 U	0.0010 U	0.0010 U	0.0031	0.0010 U	0.0010 U	0.0027	0.0010 U	0.0010 U
	HS SER-RAMW02-051415	14-May-15		0.00028 J	0.0020 U	0.0010 U	0.0072	0.0010 U	0.0010 U	0.0010 U	0.0020	0.0010 U	0.0010 U	0.0020	0.0010 U	0.0010 U
	HS SER-RAMW02-080615	6-Aug-15		0.00041 J	0.0020 U	0.0010 U	0.0029	0.0010 U	0.0010 U	0.0010 U	0.0032	0.0010 U	0.0010 U	0.0025	0.0010 U	0.0010 U
RAMW03	HS SER-RAMW03-121014	10-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00042 J	0.0010 U	0.0010 U	0.00054 J	0.0010 U	0.0010 U	
	HS SER-DUP02-121014	10-Dec-14	Field Duplicate	0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00040 J	0.0010 U	0.0010 U	0.00053 J	0.0010 U	0.0010 U	
	HS SER-RAMW03-021715	17-Feb-15		0.00051 J	0.0020 U	0.0010 U	0.00062 J	0.0010 U	0.0010 U	0.00098 J	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U	
	HS SER-DUP02-021715	17-Feb-15	Field Duplicate	0.00040 J	0.0020 U	0.0010 U	0.00064 J	0.0010 U	0.0010 U	0.00091 J	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U	
	HS SER-RAMW03-051415	14-May-15		0.00043 J	0.0020 U	0.0010 U	0.00034 J	0.0010 U	0.00085 J	0.0010 U	0.00044 J	0.0010 U	0.0010 U	0.0010	0.0010 U	0.0010 U
	HS SER-DUP02-051415	14-May-15	Field Duplicate	0.00036 J	0.0020 U	0.0010 U	0.00035 J	0.0010 U	0.0010 U	0.00043 J	0.0010 U	0.0010 U	0.0010	0.0010 U	0.0010 U	
	HS SER-RAMW03-080615	6-Aug-15		0.00043 J	0.0020 U	0.0010 U	0.00039 J	0.0010 U	0.0010 U	0.00043 J	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U	
	HS SER-DUP02-080615	6-Aug-15	Field Duplicate	0.00044 J	0.0020 U	0.0010 U	0.00041 J	0.0010 U	0.0010 U	0.00041 J	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U	
RAMW04	HS SER-RAMW04-121014	10-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00034 J	0.0010 U	0.0010 U	0.00040 J	0.0010 U	0.0010 U	
	HS SER-RAMW04-021715	17-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.00046 J	0.0010 U	0.0010 U	0.00045 J	0.0010 U	0.0010 U	0.00042 J	0.0010 U	0.0010 U	
	HS SER-RAMW04-051415	14-May-15		0.0010 U	0.0020 U	0.0010 U	0.00027 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0004 J	0.0010 U	0.0010 U	
	HS SER-RAMW04-080715	7-Aug-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00039 J	0.0010 U	0.0010 U	0.00043 J	0.0010 U	0.0010 U	
RAMW05	HS SER-RAMW05-121014	10-Dec-14		0.0035	0.0020 U	0.0022	0.00097 J	0.0010 U	0.0036	0.0010 U	0.0547	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-RAMW05-021715	17-Feb-15		0.0043	0.0020 U	0.0025	0.0034	0.0010 U	0.0097	0.0010 U	0.161	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-RAMW05-051415	14-May-15		0.0014	0.0020 U	0.00095 J	0.00046 J	0.0010 U	0.0012	0.0010 U	0.0222	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-RAMW05-080715	7-Aug-15		0.0010 U	0.0020 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.00052 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
RAMW06	HS SER-RAMW06-120914	9-Dec-14		0.0010	0.0020 U	0.0087 ^a	0.0047	0.0010 U	0.0067	0.0010 U	0.1790	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-RAMW06-021815	18-Feb-15		0.0033	0.0020 U	0.0074 ^a	0.0039	0.0010 U	0.0217	0.0010 U	0.387 ^a	0.0010 U	0.0010 U	0.00043 J	0.0010 U	0.0010 U
	HS SER-RAMW06-051315	13-May-15		0.0019	0.0020 U	0.0060	0.0032	0.0010 U	0.0032	0.0010 U	0.145	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-RAMW06-080715	7-Aug-15		0.00061 J	0.0020 U	0.0056	0.0026	0.0010 U	0.0068	0.0010 U	0.0596	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
RAMW07	HS SER-RAMW07-120914	9-Dec-14		0.0043	0.0020 U	0.0056	0.0027	0.0010 U	0.0025	0.0010 U	0.1300	0.0010 U	0.0010 U	0.0011	0.0010 U	0.0010 U
	HS SER-RAMW07-021815	18-Feb-15		0.0029	0.0020 U	0.0040	0.0091	0.0010 U	0.0053	0.0010 U	0.149	0.0010 U	0.0010 U	0.0012	0.0010 U	0.0010 U
	HS SER-RAMW07-051315	13-May-15		0.0016 J	0.010 U	0.0516 ^a	0.0584	0.0050 U	0.0705 ^a	0.0050 U	1.040 ^a	0.0050 U	0.0041 J	0.0032 J	0.0050 U	0.0050 U

Table 3
Fourth Quarter 2014 to Third Quarter 2015 Groundwater Analytical Results - Performance Wells
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

				Trichloroethene (TCE)	Methylene Chloride (Dichloromethane)	1,1-Dichloroethene	1,1-Dichloroethane	1,2-Dichloroethane	cis-1,2-Dichloroethene	trans-1,2-Dichloroethene	1,1,1-Trichloroethane	1,1,2-Trichloroethane	Ethylbenzene	Tetrachloroethene (PCE)	Toluene	Vinyl chloride	
Preliminary Remediation Goals (PRG) ^A				0.005 ^c	0.005 ^c	0.007 ^{b,c}	0.7 ^A	0.005 ^c	0.07 ^c	0.1 ^c	0.2 ^{b,c}	0.005 ^c	0.7 ^c	0.005 ^c	1.0 ^c	0.002 ^c	
Well	Sample ID	Sample Date	Sample Type	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	
RAMW08	HS SER-RAMW08-120914	9-Dec-14		0.0010 U	0.0020 U	0.0010 U	0.0014	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	
	HS SER-RAMW08-021815	18-Feb-15		0.0010 U	0.0020 U	0.0010 U	0.0018	0.0010 U	0.0010 U	0.0010 U	0.00039 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-RAMW08-051315	13-May-15		0.0010 U	0.0020 U	0.0010 U	0.0009 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U
	HS SER-RAMW08-080615	6-Aug-15		0.0010 U	0.0020 U	0.0010 U	0.00025 J	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U	0.0010 U

Notes:

- PRG Preliminary Remediation Goals (PRGs) from the Record of Decision (ROD)
 A Class 1 - Groundwater Remediation Objectives
6.5^A Concentration exceeds the indicated standard at specified well; however, compliance with the standard is only applicable to GMZ wells.
15.2 Concentration was detected but did not exceed applicable standards.
0.50 U Laboratory estimated quantitation limit exceeded standard.
 0.03 U The analyte was not detected above the laboratory estimated quantitation limit.
 mg/L milligrams per liter
 n/v No standard/guideline value.
 - Parameter not analyzed / not available.
- b,c Oral Reference Dose and/or Reference Concentration under review by USEPA. Listed values subject to change.
 Groundwater Quality Standard for this chemical pursuant to 35 Ill.Adm.Code 620.410 for Class I Groundwater or 35 Ill.Adm.Code 620.420 for Class II Groundwater.
 c Value listed is also the Groundwater Quality Standard for this chemical pursuant to 35 Ill.Adm.Code 620.410 for Class I Groundwater or 35 Ill.Adm.Code 620.420 for Class II Groundwater.
 B The analyte was detected in the method, field and/or trip blank.
 J Indicates estimated value.
 NJ The analysis indicates the presence of an analyte that has been "tentatively identified" and the associated numerical value represents its approximate concentration.

considered during evaluation of this data.

Table 4.1
Call 1 - Phase 1 SVE System Effluent Data
December 2009 - September 2015

UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 1 SVE EFFLUENT										CELL 2 SVE EFFLUENT									
Date	Sample Type	SVE Run Time (hr)	Cell 1 Run Time (hr)	SVE Flow Rate (gpm)	1,1,1-Trichloroethane	1,1,2-Trichloroethane	1,1-Dichloroethane	1,2-Dichloroethane	cis-1,2-Dichloroethane	trans-1,2-Dichloroethane	Trichloroethene	Vinyl chloride	Methylene Chloride						
				Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)		
12/1/02/09		150	53	140	13000	3.7E-02	140 U	0.00E+00	45000	6.0E-02	140 U	0.00E+00	940	3.9E-03	280	7.4E-04	8100	1.10E-02	
12/22/2009		372	124	140	860	2.6E-03	28 U	0.00E+00	11000	2.3E-02	26 U	0.00E+00	300	1.4E-03	41	1.1E-04	470	6.3E-04	
2/24/2010		1803	631	150	640	1.9E-03	6.0 U	0.00E+00	1800	4.3E-03	6.0 U	0.00E+00	150	5.7E-04	24	7.3E-05	4,80E-05	8.0 U	
3/15/2010		2345	782	140	1100	3.1E-03	8.4 U	0.00E+00	2800	6.0E-03	8.4 U	0.00E+00	37	7.7E-03	140	3.1E-03	570	9.25E-04	
4/14/2010		2804	935	150	1400	4.3E-03	12 U	0.00E+00	4100	9.4E-03	12 U	0.00E+00	31	9.0E-05	140	3.1E-03	86	3.05E-03	
5/13/2010		3405	1105	140	580	1.7E-03	7.0 U	0.00E+00	280	5.5E-03	7.0 U	0.00E+00	300	1.0E-03	32	9.1E-05	10	1.3E-05	
6/2/2010		4430	1477	108	710	1.5E-03	8.6 U	0.00E+00	2800	4.3E-03	8.6 U	0.00E+00	200	8.05E-04	30	6.0E-05	8.8 U	0.00E+00	
7/21/2010		5058	1898	140	480	1.3E-03	7.0 U	0.00E+00	2800	5.5E-03	7.0 U	0.00E+00	10	2.55E-03	42	1.2E-04	7.0 U	0.00E+00	
8/23/2010		57B4	1828	0	370	0.00E+00	8.2 U	0.00E+00	2400	0.00E+00	8.2 U	0.00E+00	500	0.00E+00	82	0.00E+00	8.2 U	0.00E+00	
9/23/2010		6523	2174	145	480	1.4E-03	7.2 U	0.00E+00	2000	4.4E-03	7.2 U	0.00E+00	250	5.4E-04	7.2 U	0.00E+00	7.2 U	0.00E+00	
10/22/2010		7219	2405	140	380	1.3E-03	5.0 U	0.00E+00	1800	3.4E-03	5.0 U	0.00E+00	240	8.6E-04	21	5.0E-05	5.0 U	0.00E+00	
10/22/2010	Dup	7219	2405	140	2800	1.2E-03	4.3 U	0.00E+00	980	3.0E-03	10 U	0.00E+00	120	1.0E-04	49	1.0E-04	10 U	0.00E+00	
11/5/2010		7794	2568	140	420	1.2E-03	4.3 U	0.00E+00	1700	3.0E-03	4.3 U	0.00E+00	140	2.0E-04	43	5.0E-05	43	0.00E+00	
12/22/2010		8508	2777	150	600	1.8E-03	4.2 U	0.00E+00	1800	3.6E-03	4.2 U	0.00E+00	8.5	1.9E-05	510	1.1E-03	42	1.00E+00	
12/22/2010		8202	2975	170	380	1.2E-04	2.3 U	0.00E+00	730	1.0E-03	2.3 U	0.00E+00	28	7.15E-06	63	1.6E-04	52	0.00E+00	
1/2/2011		10771	1857	165	280	9.5E-04	4.0 U	0.00E+00	1800	4.0E-03	4.0 U	0.00E+00	45	1.2E-05	120	2.9E-04	45	1.00E+00	
2/22/2011		10673	3263	165	200	8.8E-04	6.3 U	0.00E+00	1900	4.8E-03	6.3 U	0.00E+00	32	1.3E-04	63	8.3U	0.00E+00	4.0 U	
3/18/2011		11241	3408	160	180	5.9E-04	4.5 U	0.00E+00	1700	4.1E-03	4.5 U	0.00E+00	43	1.7E-04	68	4.5E-05	4.5 U	0.00E+00	
4/15/2011		3408	1741	160	110	3.8E-04	4.3 U	0.00E+00	1100	2.4E-04	4.3 U	0.00E+00	140	5.0E-04	16	3.3E-05	4.3 U	0.00E+00	
5/18/2011		12061	3065	160	110	3.8E-04	4.3 U	0.00E+00	1100	2.4E-04	4.3 U	0.00E+00	85	2.9E-04	11	3.3E-05	4.3 U	0.00E+00	
6/19/2011		12722	3839	170	150	5.2E-04	2.3 U	0.00E+00	730	1.9E-03	2.3 U	0.00E+00	2.8	7.15E-06	63	1.6E-04	2.3 U	0.00E+00	
7/15/2011		13417	4472	170	140	4.9E-04	1.2 U	0.00E+00	380	1.0E-03	1.2 U	0.00E+00	2.2	5.6E-08	47	1.2E-04	1.2 U	0.00E+00	
8/22/2011		14324	4775	170	150	5.2E-04	1.1 U	0.00E+00	210	5.4E-05	1.1 U	0.00E+00	2.1	5.3E-05	36	2.0E-06	1.1 U	0.00E+00	
9/15/2011		14905	4988	170	130	4.7E-04	1.1 U	0.00E+00	130	3.0E-04	1.1 U	0.00E+00	15	3.8E-08	40	1.0E-04	1.1 U	0.00E+00	
10/14/2011		15586	5199	160	65	2.1E-04	0.74 U	0.00E+00	100	2.45E-04	0.74 U	0.00E+00	14	3.37E-06	43	1.03E-04	0.86	2.07E-06	
11/2/2011		1650	5503	170	170	4.9B	1.7E-04	0.00E+00	68	0.00E+00	0.14 U	0.00E+00	2.9	7.15E-05	61	2.87E-04	7.3	2.53E-06	
12/14/2011		17010	5674	170	53	1.8E-04	0.78 U	0.00E+00	45	1.17E-04	0.78 U	0.00E+00	18	4.60E-05	78 U	0.00E+00	5.8	1.05E-06	
1/18/2012		17923	5974	170	51	1.7E-04	0.79 U	0.00E+00	41	1.07E-04	0.79 U	0.00E+00	1	2.50E-06	12	3.07E-06	170	7.9U	
2/15/2012		18506	6189	170	48	1.6E-04	0.78 U	0.00E+00	30	7.82E-05	0.78 U	0.00E+00	37	1.82E-04	5	1.4E-05	1.4E-04	4.2E-06	
3/15/2012		18262	6421	160	38	1.3E-04	0.71 U	0.00E+00	34	8.67E-05	0.78 U	0.00E+00	32	1.35E-05	71 U	0.00E+00	3.9	1.40E-04	
4/19/2012		20748	6919	160	51	1.6E-04	0.76 U	0.00E+00	38	8.84E-05	0.78 U	0.00E+00	58	2.98E-04	61	1.89E-05	0.76 U	0.00E+00	
Pulse-off period: June 1, 2012 to August 14, 2012										Pulse-off period: June 1, 2012 to August 14, 2012									
8/14/2012		21282	7004	160	120	3.9E-04	1.3 U	0.00E+00	51	1.2E-04	1.3 U	0.00E+00	2	4.81E-05	13 U	0.00E+00	13 U	0.00E+00	
9/17/2012		21062	7317																

Table 4.1
Cell 1 - Phase 1 SVE System Effluent Data
December 2008 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 1 SVE EFFLUENT																	
Date	Sample Type	SVE Run Time (hr)	Cell 1 Run Time (hr)	SVE Flow Rate (gpm)	Conc (ppbv) Removed Rate (lb/hr)	Mass Conc (ppbv)	Chloroethane	Benzene	Toluene	Ethylbenzene	m,p-Xylene	o-Xylenes	Aromate	Methyl Ethyl Ketone (MEK)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal Rate (lb)	
12/1/2009		159	53	140	140U 0.00E+00	140U 0.00E+00	17000	2.38E-02	140U 0.00E+00	250 5.78E-04	4.15E-03	470 1.08E-03	3800 4.78E-03	140U 0.00E+00	2.28E-01	11.81	
12/2/2009		372	124	140	26U 0.00E+00	26U 0.00E+00	1700	2.38E-03	28U 0.00E+00	32 8.60E-05	28U 0.00E+00	470 0.00E+00	100U 0.00E+00	28U 0.00E+00	4.67E-02	15.23	
2/2/2010		1863	631	150	6.0U 0.00E+00	6.0U 0.00E+00	130	1.05E-04	19 3.45E-05	6.0U 0.00E+00	6.0U 0.00E+00	370 1.32E-04	370 0.00E+00	370 0.00E+00	0.95E-03	20.06	
3/15/2010		2345	782	140	8.4U 0.00E+00	8.4U 0.00E+00	170	2.38E-04	8.4U 0.00E+00	8.4U 0.00E+00	34U 8.4U	34U 0.00E+00	34U 0.00E+00	1.30E-02	22.02		
4/1/2010		2804	935	150	12U 0.00E+00	12U 0.00E+00	320	4.80E-04	14 2.54E-05	12U 0.00E+00	12U 0.00E+00	50U 12U	50U 0.00E+00	50U 0.00E+00	2.10E-02	28.22	
5/15/2010		3495	1185	140	7.0U 0.00E+00	7.0U 0.00E+00	100	1.40E-04	12 2.03E-05	7.0U 0.00E+00	7.0U 0.00E+00	28U 7.0U	28U 0.00E+00	28U 0.00E+00	1.10E-02	27.75	
6/2/2010		4430	1477	108	8.8U 0.00E+00	8.8U 0.00E+00	80	9.40E-05	10 1.31E-05	8.8U 0.00E+00	8.8U 0.00E+00	34J 8.8U	34J 0.00E+00	34J 0.00E+00	7.88E-03	30.20	
7/2/2010		5058	1886	140	7.0U 0.00E+00	7.0U 0.00E+00	24	8.40E-05	7.0U 0.00E+00	7.0U 0.00E+00	7.0U 7.0U	7.0U 0.00E+00	7.0U 0.00E+00	1.11E-02	32.52		
8/2/2010		5784	1928	0	8.2U 0.00E+00	8.2U 0.00E+00	82U 8.2U	0.00E+00	82U 0.00E+00	82U 0.00E+00	53 82U	53 0.00E+00	53 0.00E+00	0.00E+00	32.52		
9/2/2010		6623	2174	145	7.2U 0.00E+00	7.2U 0.00E+00	72U 7.2U	0.00E+00	72U 0.00E+00	72U 0.00E+00	72U 7.2U	72U 0.00E+00	72U 0.00E+00	7.98E-03	34.49		
10/2/2010	Dup	7219	2408	140	5.0U 0.00E+00	5.0U 0.00E+00	11 1.0U	0.00E+00	10U 0.00E+00	10U 0.00E+00	50U 5.0U	50U 0.00E+00	50U 0.00E+00	5.91E-03	35.86		
10/2/2010		7219	2408	140	10U 0.00E+00	10U 0.00E+00	12 1.0U	0.00E+00	12 1.0U	0.00E+00	43U 4.3U	43U 0.00E+00	43U 0.00E+00	4.31E-03	36.96		
11/15/2010		7794	2598	140	4.3U 0.00E+00	4.3U 0.00E+00	10 1.0U	0.00E+00	10 1.0U	0.00E+00	42U 4.2U	42U 0.00E+00	42U 0.00E+00	2.18E-05	38.22		
12/2/2010		8508	2777	150	4.2U 0.00E+00	4.2U 0.00E+00	92U 5.3	0.00E+00	92U 0.00E+00	92U 0.00E+00	52U 5.2U	52U 0.00E+00	52U 0.00E+00	6.28E-03	39.47		
1/2/2011		9302	2975	170	5.2U 0.00E+00	5.2U 0.00E+00	18U 4.0U	0.00E+00	18U 0.00E+00	18U 0.00E+00	40U 4.0U	40U 0.00E+00	40U 0.00E+00	4.00E-03	40.53		
2/2/2011		10071	3167	165	4.0U 0.00E+00	4.0U 0.00E+00	25U 6.3U	0.00E+00	25U 6.3U	0.00E+00	63U 6.3U	63U 0.00E+00	63U 0.00E+00	5.67E-03	41.53		
3/1/2011		10573	3283	165	6.3U 0.00E+00	6.3U 0.00E+00	18U 4.5U	0.00E+00	18U 4.5U	0.00E+00	45U 4.5U	45U 0.00E+00	45U 0.00E+00	4.52E-03	42.15		
4/15/2011		11241	3460	160	4.3U 0.00E+00	4.3U 0.00E+00	17U 4.3U	0.00E+00	17U 4.3U	0.00E+00	43U 4.3U	43U 0.00E+00	43U 0.00E+00	3.52E-03	42.87		
5/16/2011		12081	3665	160	4.3U 0.00E+00	4.3U 0.00E+00	92U 2.3U	0.00E+00	92U 2.3U	0.00E+00	23U 2.3U	23U 0.00E+00	23U 0.00E+00	3.12E-03	43.39		
6/16/2011		12722	3830	170	2.3U 0.00E+00	2.3U 0.00E+00	30U 3.0U	0.00E+00	30U 3.0U	0.00E+00	23U 2.3U	23U 0.00E+00	23U 0.00E+00	9.21U	0.00E+00		
7/16/2011		13417	4472	170	1.2U 0.00E+00	1.2U 0.00E+00	12U 1.2U	0.00E+00	12U 1.2U	0.00E+00	12U 1.2U	0.00E+00	12U 1.2U	0.00E+00	2.44E-03	44.96	
8/2/2011		14324	4775	170	1.1U 0.00E+00	1.1U 0.00E+00	11U 1.1U	0.00E+00	11U 1.1U	0.00E+00	11U 1.1U	0.00E+00	11U 1.1U	0.00E+00	45.59	45.59	
9/15/2011		14905	4988	170	1.1U 0.00E+00	1.1U 0.00E+00	11U 1.1U	0.00E+00	11U 1.1U	0.00E+00	18U 1.1U	18U 0.00E+00	18U 0.00E+00	18U 0.00E+00	5.97E-03	45.93	
10/1/2011		15588	5199	160	0.74U 0.00E+00	0.74U 0.00E+00	30U 3.0U	0.00E+00	30U 3.0U	0.00E+00	17U 1.8	17U 0.00E+00	17U 0.00E+00	17U 0.00E+00	4.21E-03	46.20	
11/2/2011		16510	5503	170	0.74U 0.00E+00	0.74U 0.00E+00	74U 7.4U	0.00E+00	74U 7.4U	0.00E+00	74U 7.4U	74U 0.00E+00	74U 0.00E+00	14 5.68E-04	46.37		
12/1/2011		17010	5674	170	0.78U 0.00E+00	0.78U 0.00E+00	31U 3.1U	0.00E+00	31U 3.1U	0.00E+00	78U 7.8U	78U 0.00E+00	78U 0.00E+00	7.8U 0.00E+00	4.80E-06	46.48	
1/16/2012		17923	18580	170	0.78U 0.00E+00	0.78U 0.00E+00	32U 3.2U	0.00E+00	32U 3.2U	0.00E+00	78U 7.8U	78U 0.00E+00	78U 0.00E+00	20 3.08E-05	3.1U 0.00E+00	5.01E-04	
2/18/2012		21518	6189	170	0.78U 0.00E+00	0.78U 0.00E+00	12U 1.2U	0.00E+00	12U 1.2U	0.00E+00	78U 7.8U	78U 0.00E+00	78U 0.00E+00	2.8U 0.00E+00	4.02E-04	46.84	
3/15/2012		18202	6421	170	0.71U 0.00E+00	0.71U 0.00E+00	30U 3.0U	0.00E+00	30U 3.0U	0.00E+00	78U 7.8U	78U 0.00E+00	78U 0.00E+00	5.50E-04	46.86		
4/18/2012		21092	20748	160	0.76U 0.00E+00	0.76U 0.00E+00	30U 3.0U	0.00E+00	30U 3.0U	0.00E+00	78U 7.8U	78U 0.00E+00	78U 0.00E+00	6.11E-04	47.12		
Pulse-off period June 1, 2012 to August 14, 2012																	
8/17/2012		21282	7094	160	1.3U 0.00E+00	1.3U 0.00E+00	13U 1.3U	0.00E+00	13U 1.3U	0.00E+00	13U 1.3U	0.00E+00	13U 1.3U	0.00E+00	5.2U 0.00E+00	2.03E-03	47.48

Table 4.2
Cell 2 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 2 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 2 Run Time (hr)	SVE Flow Rate (scfm)	1,1,1-Trichloroethane	1,1,2-Trichloroethane	1,1-Dichloroethane	1,2-Dichloroethane	1,1-Dichloroethene	cis-1,2-Dichloroethene	trans-1,2-Dichloroethene	Tetrachloroethene
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
12/11/2009		178	59	150	40000	1.24E-01	86 U	0.00E+00	21000	4.83E-02	86 U	0.00E+00
12/15/2009		205	68	140	27000	7.82E-02	110 U	0.00E+00	14000	3.01E-02	110 U	0.00E+00
12/29/2009		539	180	140	24000	6.95E-02	100 U	0.00E+00	9100	1.95E-02	100 U	0.00E+00
1/13/2010		903	301	150	9100	2.82E-02	35 U	0.00E+00	3700	8.52E-03	35 U	0.00E+00
1/27/2010		1224	408	150	13000	4.03E-02	40 U	0.00E+00	4300	9.90E-03	40 U	0.00E+00
1/27/2010	Dup	1224	408	150	14000	4.34E-02	40 U	0.00E+00	4800	1.10E-02	40 U	0.00E+00
2/24/2010		1893	631	150	8000	2.48E-02	22 U	0.00E+00	3000	6.90E-03	22 U	0.00E+00
3/15/2010		2345	782	140	17000	4.92E-02	48 U	0.00E+00	8000	1.72E-02	48 U	0.00E+00
4/14/2010		2804	935	150	8400	2.61E-02	23 U	0.00E+00	2200	5.06E-03	23 U	0.00E+00
5/13/2010		3495	1165	140	8000	2.32E-02	11 U	0.00E+00	3100	6.66E-03	11 U	0.00E+00
6/21/2010		4430	1477	108	5800	1.30E-02	23 U	0.00E+00	3000 J	4.97E-03	23 U	0.00E+00
7/21/2010		5058	1686	140	4500	1.30E-02	14 U	0.00E+00	1600	3.44E-03	14 U	0.00E+00
8/23/2010		5784	1928	0	7100	0.00E+00	20 U	0.00E+00	2700	0.00E+00	20 U	0.00E+00
9/23/2010		6523	2174	145	4300	1.29E-02	12 U	0.00E+00	1600	3.56E-03	12 U	0.00E+00
10/22/2010		7219	2406	140	2500	7.24E-03	10 U	0.00E+00	890	1.91E-03	10 U	0.00E+00
11/15/2010		7794	2598	140	3200	9.27E-03	11 U	0.00E+00	1100	2.36E-03	11 U	0.00E+00
12/22/2010		8508	2955	150	4000	1.24E-02	14 U	0.00E+00	1500	3.45E-03	14 U	0.00E+00
1/24/2011		9302	3352	170	780	2.74E-03	2.7 U	0.00E+00	800	2.09E-03	2.7 U	0.00E+00
2/25/2011		10071	3737	165	1500	5.12E-03	4.0 U	0.00E+00	1100	2.78E-03	4.0 U	0.00E+00
3/18/2011		10573	3988	165	370	1.26E-03	1.0 U	0.00E+00	160	4.05E-04	1.0 U	0.00E+00
4/15/2011		11241	4322	160	300 J,B	9.93E-04	1.0 U	0.00E+00	95	2.33E-04	1.0 U	0.00E+00
5/19/2011		12061	4732	160	93	3.08E-04	1.1 U	0.00E+00	39	9.57E-05	1.1 U	0.00E+00
6/16/2011		12722	5062	170	99	3.48E-04	1.2 U	0.00E+00	48	1.25E-04	1.2 U	0.00E+00
7/15/2011		13417	4472	170	77	2.71E-04	1.2 U	0.00E+00	25	6.52E-05	1.2 U	0.00E+00
8/22/2011		14324	4775	170	78	2.74E-04	1.2 U	0.00E+00	31	8.09E-05	1.2 U	0.00E+00
9/15/2011		14905	4968	170	69	2.43E-04	1.1 U	0.00E+00	20	5.22E-05	1.1 U	0.00E+00
10/14/2011		15598	5199	160	43	1.42E-04	0.82 U	0.00E+00	12	2.95E-05	0.82 U	0.00E+00
11/21/2011		16510	5503	170	28 J,B	9.85E-05	1.6 U	0.00E+00	7.7	2.01E-05	1.6 U	0.00E+00
12/14/2011		17010	5670	170	26	9.14E-05	0.76 U	0.00E+00	5.2	1.36E-05	0.76 U	0.00E+00
1/19/2012		17923	5974	170	25	8.79E-05	0.74 U	0.00E+00	6.9	1.80E-05	0.74 U	0.00E+00
2/15/2012		18566	6189	170	31	1.09E-04	0.73 U	0.00E+00	7.6	1.98E-05	0.73 U	0.00E+00
3/15/2012		19262	6421	170	29 J,B	1.02E-04	0.71 U	0.00E+00	8.1	2.11E-05	0.71 U	0.00E+00
4/19/2012		20102	6701	160	31	1.03E-04	0.76 U	0.00E+00	11	2.70E-05	0.76 U	0.00E+00
5/16/2012		20748	6916	160	28	9.27E-05	0.78 U	0.00E+00	11	2.70E-05	0.78 U	0.00E+00



Table 4.2
Cell 2 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 2 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 2 Run Time (hr)	SVE Flow Rate (scfm)	Trichloroethene		Vinyl chloride		Methylene Chloride		Carbon Tetrachloride		Chloroform		Chloroethane		Benzene		Toluene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
12/11/2009		178	59	150	330	1.01E-03	4400	6.40E-03	86 U	0.00E+00	86 U	0.00E+00	86 U	0.00E+00	210	3.15E-04	86 U	0.00E+00	200	4.29E-04
12/15/2009		205	68	140	240	6.84E-04	3500	4.75E-03	110 U	0.00E+00	110 U	0.00E+00	110 U	0.00E+00	370	5.18E-04	110 U	0.00E+00	140	2.80E-04
12/29/2009		539	180	140	240	6.84E-04	1500	2.03E-03	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	120	1.68E-04	100 U	0.00E+00	100 U	0.00E+00
1/13/2010		903	301	150	130	3.97E-04	250	3.63E-04	35 U	0.00E+00	35 U	0.00E+00	35 U	0.00E+00	170	2.55E-04	35 U	0.00E+00	35 U	0.00E+00
1/27/2010		1224	408	150	150	4.58E-04	200	2.91E-04	40 U	0.00E+00	40 U	0.00E+00	40 U	0.00E+00	120	1.80E-04	40 U	0.00E+00	40 U	0.00E+00
1/27/2010	Dup	1224	408	150	180	5.50E-04	240	3.49E-04	40 U	0.00E+00	40 U	0.00E+00	40 U	0.00E+00	130	1.95E-04	40 U	0.00E+00	40 U	0.00E+00
2/24/2010		1893	631	150	98	2.99E-04	73	1.06E-04	22 U	0.00E+00	22 U	0.00E+00	22 U	0.00E+00	38	5.70E-05	22 U	0.00E+00	22 U	0.00E+00
3/15/2010		2345	782	140	210	5.99E-04	62	8.41E-05	48 U	0.00E+00	48 U	0.00E+00	48 U	0.00E+00	180	2.52E-04	48 U	0.00E+00	48 U	0.00E+00
4/14/2010		2804	935	150	190	5.81E-04	69	1.00E-04	23 U	0.00E+00	23 U	0.00E+00	23 U	0.00E+00	23 U	0.00E+00	23 U	0.00E+00	23 U	0.00E+00
5/13/2010		3495	1165	140	78	2.22E-04	42	5.70E-05	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	20	2.80E-05	11 U	0.00E+00	11 U	0.00E+00
6/21/2010		4430	1477	108	88	1.94E-04	23 U	0.00E+00	23 U	0.00E+00	23 U	0.00E+00	23 U	0.00E+00	33 J	3.56E-05	23 U	0.00E+00	23 U	0.00E+00
7/21/2010		5058	1686	140	80	2.28E-04	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00
8/23/2010		5784	1928	0	150	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	21	0.00E+00	20 U	0.00E+00	20 U	0.00E+00
9/23/2010		6523	2174	145	74	2.19E-04	12	1.69E-05	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00
10/22/2010		7219	2406	140	42	1.20E-04	10 U	0.00E+00	10 U	0.00E+00	10 U	0.00E+00	10 U	0.00E+00	10 U	0.00E+00	10 U	0.00E+00	10 U	0.00E+00
11/15/2010		7794	2598	140	35	9.98E-05	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00
12/22/2010		8508	2955	150	27	8.25E-05	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00
1/24/2011		9302	3352	170	9	3.12E-05	2.7 U	0.00E+00	2.7 U	0.00E+00	2.7 U	0.00E+00	2.7 U	0.00E+00	2.7 U	0.00E+00	2.7 U	0.00E+00	3.1	7.53E-06
2/25/2011		10071	3737	165	15	5.04E-05	4.0 U	0.00E+00	4.0 U	0.00E+00	4.0 U	0.00E+00	4.0 U	0.00E+00	16 U	0.00E+00	4.0 U	0.00E+00	4.0 U	0.00E+00
3/18/2011		10573	3988	165	7.3	2.45E-05	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	4.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00
4/15/2011		11241	4322	160	8.5	2.77E-05	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	4.1 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00
5/19/2011		12061	4732	160	11	3.59E-05	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.5 U	0.00E+00	8.1	1.57E-05	1.1 U	0.00E+00
6/16/2011		12722	5062	170	15	5.19E-05	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.9	3.91E-06	1.2 U	0.00E+00
7/15/2011		13417	4472	170	21	7.27E-05	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.6 U	0.00E+00	5.3	1.09E-05	1.2 U	0.00E+00
8/22/2011		14324	4775	170	22	7.62E-05	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.6	3.29E-06	1.2 U	0.00E+00
9/15/2011		14905	4968	170	18	6.23E-05	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.5 U	0.00E+00	4.6	9.47E-06	1.1 U	0.00E+00
10/14/2011		15598	5199	160	9.1	2.97E-05	0.82 U	0.00E+00	0.82 U	0.00E+00	0.82 U	0.00E+00	0.82 U	0.00E+00	3.3 U	0.00E+00	0.82 U	0.00E+00	0.82 U	0.00E+00
11/21/2011		16510	5503	170	5.1	1.77E-05	1.6 U	0.00E+00	1.6 UJ	0.00E+00	1.6 U	0.00E+00	1.6 U	0.00E+00	6.4 U	0.00E+00	1.6 U	0.00E+00	1.6 U	0.00E+00
12/14/2011		17010	5670	170	3.4	1.18E-05	0.76 U	0.00E+00	7.6 U	0.00E+00	0.76 U	0.00E+00	0.76 U	0.00E+00	3.0 U	0.00E+00	0.78	1.61E-06	0.76 U	0.00E+00
1/19/2012		17923	5974	170	2.9	1.00E-05	0.74 U	0.00E+00	1.9	4.25E-06	0.74 U	0.00E+00								

Table 4.2
Cell 2 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 2 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 2 Run Time (hr)	SVE Flow Rate (scfm)	Ethylbenzene		m&p-Xylenes		o-Xylenes		Acetone		Methyl Ethyl Ketone (MEK)		Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)		
12/11/2009		178	59	150	86 U	0.00E+00	240	5.93E-04	110	2.72E-04	340 U	0.00E+00	86 U	0.00E+00	2.54E-01	15.05
12/15/2009		205	68	140	110 U	0.00E+00	230	5.30E-04	110 U	0.00E+00	430 U	0.00E+00	110 U	0.00E+00	1.59E-01	16.48
12/29/2009		539	180	140	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	420 U	0.00E+00	100 U	0.00E+00	1.19E-01	29.76
1/13/2010		903	301	150	35 U	0.00E+00	35 U	0.00E+00	35 U	0.00E+00	140 U	0.00E+00	35 U	0.00E+00	4.93E-02	35.75
1/27/2010		1224	408	150	40 U	0.00E+00	40 U	0.00E+00	40 U	0.00E+00	160 U	0.00E+00	40 U	0.00E+00	6.47E-02	42.68
1/27/2010	Dup	1224	408	150	40 U	0.00E+00	40 U	0.00E+00	40 U	0.00E+00	160 U	0.00E+00	40 U	0.00E+00	7.06E-02	43.31
2/24/2010		1893	631	150	22 U	0.00E+00	22 U	0.00E+00	22 U	0.00E+00	87 U	0.00E+00	22 U	0.00E+00	3.93E-02	51.44
3/15/2010		2345	782	140	48 U	0.00E+00	48 U	0.00E+00	48 U	0.00E+00	190 U	0.00E+00	48 U	0.00E+00	8.60E-02	64.40
4/14/2010		2804	935	150	23 U	0.00E+00	23 U	0.00E+00	23 U	0.00E+00	92 U	0.00E+00	23 U	0.00E+00	4.24E-02	70.89
5/13/2010		3495	1165	140	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	43 U	0.00E+00	11 U	0.00E+00	3.84E-02	79.74
6/21/2010		4430	1477	108	23 U	0.00E+00	23 U	0.00E+00	23 U	0.00E+00	92 U	0.00E+00	23 U	0.00E+00	2.30E-02	86.90
7/21/2010		5058	1686	140	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	58 U	0.00E+00	14 U	0.00E+00	2.07E-02	91.24
8/23/2010		5784	1928	0	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	81 U	0.00E+00	20 U	0.00E+00	0.00E+00	91.24
9/23/2010		6523	2174	145	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	47 U	0.00E+00	12 U	0.00E+00	2.04E-02	96.27
10/22/2010		7219	2406	140	10 U	0.00E+00	10 U	0.00E+00	10 U	0.00E+00	42 U	0.00E+00	10 U	0.00E+00	1.11E-02	98.85
11/15/2010		7794	2598	140	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	44 U	0.00E+00	11 U	0.00E+00	1.34E-02	101.41
12/22/2010		8508	2955	150	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	56 U	0.00E+00	14 U	0.00E+00	1.84E-02	107.99
1/24/2011		9302	3352	170	2.7 U	0.00E+00	2.7 U	0.00E+00	2.7 U	0.00E+00	11 U	0.00E+00	11	2.09E-05	6.06E-03	110.39
2/25/2011		10071	3737	165	4.0 U	0.00E+00	4.0 U	0.00E+00	4.0 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	9.59E-03	114.08
3/18/2011		10573	3988	165	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	15	2.23E-05	4.0 U	0.00E+00	1.98E-03	114.57
4/15/2011		11241	4322	160	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	8.2 J,B	1.18E-05	4.1 U	0.00E+00	1.48E-03	115.07
5/19/2011		12061	4732	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11	1.58E-05	4.5 U	0.00E+00	5.87E-04	115.31
6/16/2011		12722	5062	170	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	19	2.91E-05	4.7 U	0.00E+00	7.49E-04	115.55
7/15/2011		13417	4472	170	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	19	2.91E-05	4.6 U	0.00E+00	6.30E-04	115.18
8/22/2011		14324	4775	170	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	6.8 J,B	1.04E-05	4.7 U	0.00E+00	7.28E-04	115.40
9/15/2011		14905	4968	170	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11	1.68E-05	4.5 U	0.00E+00	5.54E-04	115.51
10/14/2011		15598	5199	160	0.82 U	0.00E+00	0.82 U	0.00E+00	0.82 U	0.00E+00	5	7.20E-06	3.3 U	0.00E+00	2.58E-04	115.57
11/21/2011		16510	5503	170	1.6 U	0.00E+00	1.6 U	0.00E+00	1.6 U	0.00E+00	6.4 U,J	0.00E+00	6.4 U	0.00E+00	1.77E-04	115.62
12/14/2011		17010	5670	170	0.76 U	0.00E+00	0.76 U	0.00E+00	0.76 U	0.00E+00	7.6 U,J	0.00E+00	3.0 U	0.00E+00	1.65E-04	115.65
1/19/2012		17923	5974	170	0.79	2.21E-06	1.5	4.20E-06	1.1	3.08E-06	14	2.14E-05	3.0 U	0.00E+00	1.80E-04	115.71
2/15/2012		18566	6189	170	0.73 U	0.00E+00	0.73 U	0.00E+00	0.73 U	0.00E+00	7.9	1.21E-05	2.9 U	0.00E+00	1.83E-04	115.74
3/15/2012		19262	6421	170	0.71 U	0.00E+00	0.71 U	0.00E+00	0.71 U	0.00E+00	8.9	1.36E-05	2.8 U	0.00E+00	1.75E-04	115.79
4/19/2012		20102	6701	160	0.76 U	0.00E+00	0.76 U	0.00E+00	0.76 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00	1.88E-04	115.84
5/16/2012		20748	6916	160	0.78 U	0.00E+00	0.78 U	0.00E+00	0.78 U	0.00E+00	3.1 U	0.00E+00	3.1 U	0.00E+00	1.94E-04	115.88

Table 4.2
Cell 2 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 2 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 2 Run Time (hr)	SVE Flow Rate (scfm)	1,1,1-Trichloroethane		1,1,2-Trichloroethane		1,1-Dichloroethane		1,2-Dichloroethane		1,1-Dichloroethene		cis-1,2-Dichloroethene		trans-1,2-Dichloroethene		Tetrachloroethene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
Pulse -off period June 1, 2012 to August 14, 2012																				
8/14/2012		21282	7094	160	370	1.22E-03	1.3 U	0.00E+00	160	3.93E-04	1.3 U	0.00E+00	4.5	1.08E-05	11	2.65E-05	1.3 U	0.00E+00	16	6.58E-05
9/17/2012		21952	7317	160	180	5.96E-04	1.1 U	0.00E+00	36	8.84E-05	1.1 U	0.00E+00	1.1 U	0.00E+00	10	2.40E-05	1.1 U	0.00E+00	20	8.23E-05
Pulse -off period September 17, 2012 to November 15, 2012																				
11/15/2012		21959	7320	160	430	1.42E-03	1.1 U	0.00E+00	140	3.44E-04	1.1 U	0.00E+00	12	2.89E-05	13	3.13E-05	1.1 U	0.00E+00	41	1.69E-04
12/14/2012		22554	7518	170	610	2.14E-03	1.9 U	0.00E+00	11	2.87E-05	1.9 U	0.00E+00	12	3.07E-05	4	1.02E-05	1.9 U	0.00E+00	5.3	2.32E-05
Pulse -off period December 14, 2012 to February 26, 2013																				
2/26/2013		22556	7518	160	1.9	6.29E-06	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
4/11/2013		23581	7723	160	140	4.63E-04	1.2 U	0.00E+00	10	2.45E-05	1.2 U	0.00E+00	4	9.62E-06	3.3	7.94E-06	1.2 U	0.00E+00	5.4	2.22E-05
Pulse -off period April 11, 2013 to May 10, 2013																				
5/10/2013		23583	7724	160	210	6.95E-04	1.1 U	0.00E+00	62	1.52E-04	1.2 U	0.00E+00	3.9	9.38E-06	5.4	1.30E-05	1.1 U	0.00E+00	6.4	2.63E-05
7/15/2013		25160	8039	160	160	5.29E-04	1.1 U	0.00E+00	20	4.91E-05	1.1 U	0.00E+00	3.7	8.90E-06	3.7	8.90E-06	1.1 U	0.00E+00	21	8.64E-05
7/15/2013	Dup	25160	8039	160	160	5.29E-04	1.2 U	0.00E+00	20	4.91E-05	1.2 U	0.00E+00	3.2	7.70E-06	3.5	8.42E-06	1.2 U	0.00E+00	17	6.99E-05
Pulse -off period July 15, 2013 to September 9, 2013																				
9/9/2013		25162	8040	160	380	1.26E-03	2.0 U	0.00E+00	110	2.70E-04	2.0 U	0.00E+00	3.4	8.18E-06	7	1.68E-05	2.0 U	0.00E+00	49	2.02E-04
11/18/2013		26825	8372	160	44	1.46E-04	1.1 U	0.00E+00	11	2.70E-05	1.1 U	0.00E+00	1.3	3.13E-06	2.3	5.53E-06	1.1 U	0.00E+00	14	5.76E-05
Pulse -off period November 18, 2013 to January 15, 2014																				
1/15/2014		28218	8651	160	160	5.29E-04	1.2 U	0.00E+00	55	1.35E-04	1.2 U	0.00E+00	3.3	7.94E-06	2.9	6.97E-06	1.2 U	0.00E+00	7.2	2.96E-05
3/14/2014		29432	8894	160	16	5.29E-05	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2	0.00E+00	1.9	4.57E-06	1.2 U	0.00E+00	1.2 U	0.00E+00
3/14/2014	Dup	29432	8894	160	19	6.29E-05	1.2 U	0.00E+00	1.6	3.93E-06	1.2 U	0.00E+00	1.2 U	0.00E+00	1.8	4.33E-06	1.2 U	0.00E+00	1.7	6.99E-06
Pulse -off period March 14, 2014 to May 15, 2014																				
5/15/2014		29914	8990	160	240	7.94E-04	1.1 U	0.00E+00	99	2.43E-04	1.1 U	0.00E+00	4.8	1.15E-05	7.8	1.88E-05	1.1 U	0.00E+00	14	5.76E-05
7/23/2014		31567	9321	160	89	2.95E-04	1.2 U	0.00E+00	20	4.91E-05	1.2 U	0.00E+00	1.8	4.33E-06	3.7	8.90E-06	1.2 U	0.00E+00	11	4.52E-05
Pulse -off period July 23, 2014 to September 16, 2014																				
9/16/2014		32432	9494	160	310	1.03E-03	2.1 U	0.00E+00	120	2.95E-04	2.1 U	0.00E+00	3.9	9.38E-06	6	1.44E-05	2.1 U	0.00E+00	19	7.82E-05
11/14/2014		33847	9777	160	42	1.39E-04	1.1 U	0.00E+00	7.8	1.91E-05	1.1 U	0.00E+00	1.1 U	0.00E+00	1.6	3.85E-06	1.1 U	0.00E+00	11	4.52E-05
Pulse -off period November 14, 2014 to January 9, 2015																				
1/9/2015		33855	9778	160	210	6.95E-04	1.2 U	0.00E+00	69	1.69E-04	1.2 U	0.00E+00	3.7	8.90E-06	3.4	8.18E-06	1.2 U	0.00E+00	8.2	3.37E-05
3/13/2015		35189	10045	160	18	5.96E-05	1.3 U	0.00E+00	5.4	1.33E-05	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	3.5	1.44E-05
Pulse -off period March 13, 2015 to May 15, 2015																				
5/15/2015		35194	10046	160	240	7.94E-04	1.2 U	0.00E+00												

Table 4.2
Cell 2 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 2 SVE EFFLUENT

Date		Sample Type	SVE Run Time (hr)	Cell 2 Run Time (hr)	SVE Flow Rate (scfm)	Trichloroethene		Vinyl chloride		Methylene Chloride		Carbon Tetrachloride		Chloroform		Chloroethane		Benzene		Toluene	
						Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
Pulse-off period	June 1, 2012 to August 14, 2012																				
8/14/2012		21282	7094	160	12	3.91E-05	1.3 U	0.00E+00	13 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	5.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	
9/17/2012		21952	7317	160	29	9.45E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.3 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	
Pulse-off period	September 17, 2012 to November 15, 2012																				
11/15/2012		21959	7320	160	26	8.47E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.5 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	
12/14/2012		22554	7518	170	17	5.89E-05	1.9 U	0.00E+00	19 U	0.00E+00	1.9 U	0.00E+00	1.9 U	0.00E+00	7.4 U	0.00E+00	1.9 U	0.00E+00	1.9 U	0.00E+00	
Pulse-off period	December 14, 2012 to February 26, 2013																				
2/26/2013		22556	7518	160	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	5 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	
4/11/2013		23581	7723	160	8	2.61E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	
Pulse-off period	April 11, 2013 to May 10, 2013																				
5/10/2013		23583	7724	160	9.5	3.10E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	18	2.88E-05	1.1 U	0.00E+00	1.1 U	0.00E+00	
7/15/2013		25160	8039	160	24	7.82E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.6 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	
7/15/2013	Dup	25160	8039	160	24	7.82E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	
Pulse-off period	July 15, 2013 to September 9, 2013																				
9/9/2013		25162	8040	160	31	1.01E-04	2.0 U	0.00E+00	20 U	0.00E+00	2.0 U	0.00E+00	2.0 U	0.00E+00	8.0 U	0.00E+00	2.0 U	0.00E+00	2.0 U	0.00E+00	
11/18/2013		26825	8372	160	8.4	2.74E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.6 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	
Pulse-off period	November 18, 2013 to January 15, 2014																				
1/15/2014		28218	8651	160	7.2	2.35E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.2 U	0.00E+00	4.9	1.12E-05	
3/14/2014		29432	8894	160	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.9 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	
3/14/2014	Dup	29432	8894	160	1.5	4.89E-06	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	5.0 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	
Pulse-off period	March 14, 2014 to May 15, 2014																				
5/15/2014		29914	8990	160	6.6	2.15E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	16	2.56E-05	1.1 U	0.00E+00	1.1 U	0.00E+00	
7/23/2014		31567	9321	160	19	6.19E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.6 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	
Pulse-off period	July 23, 2014 to September 16, 2014																				
9/16/2014		32432	9494	160	26	8.47E-05	2.1 U	0.00E+00	21 U	0.00E+00	2.1 U	0.00E+00	2.1 U	0.00E+00	8.3 U	0.00E+00	3.5	6.78E-06	2.1 U	0.00E+00	
11/14/2014		33847	9777	160	7.3	2.38E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.6 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	
Pulse-off period	November 14, 2014 to January 9, 2015																				
1/9/2015		33855	9778	160	9.3	3.03E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.8 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	
3/13/2015		35189	10045	160	3.0	9.78E-06	1.3 U	0.00E+00	13 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	5.0 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	
Pulse-off period	March 13, 2015 to May 15, 2015																				
5/15/2015		35194	10046	160	5.4	1.76E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	7.0	1.12E-05	1.2 U	0.00E+00	1.2 U	0.00E+00	
7/16/2015		36677	10343	160	18.0	5.87E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.6 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	
Pulse-off period	July 16, 2015 to September 22, 2015																				
9/22/2015		36680	10343	160	30	9.78E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.5 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	

Notes:

Mass removal rate = (flow rate in scfm)(concentration in ppmv)(60/MW) / (387*1000000)

"U" indicates non-detection at the specified reporting limit; for ND compounds, zero is used in mass removal calculations.

molecular weight (values from the U.S. National Library of Chemical Structure Data).

MW molecular weight (values from the U.S. National Library of Medicine)

SCFM = standard cubic feet per minute

SCFM standard cubic feet per minute

B The analyte was detected in the method, field and/or laboratory.

When a duplicate sample was collected, the original sample result

are used in the mass calculations.

Table 4.2
Cell 2 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 2 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 2 Run Time (hr)	SVE Flow Rate (scfm)	Ethylbenzene		m&p-Xylenes		o-Xylenes		Acetone		Methyl Ethyl Ketone (MEK)		Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)		
Pulse-off period June 1, 2012 to August 14, 2012																
8/14/2012		21282	7094	160	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	21	3.03E-05	5.3 U	0.00E+00	1.79E-03	116.20
9/17/2012		21952	7317	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.3 U	0.00E+00	8.85E-04	116.40
Pulse-off period September 17, 2012 to November 15, 2012																
11/15/2012		21959	7320	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.5 U	0.00E+00	2.08E-03	116.40
12/14/2012		22554	7518	170	1.9 U	0.00E+00	1.9 U	0.00E+00	1.9 U	0.00E+00	19 U	0.00E+00	7.4 U	0.00E+00	2.30E-03	116.86
Pulse-off period December 14, 2012 to February 26, 2013																
2/26/2013		22556	7518	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	5 U	0.00E+00	6.29E-06	116.86
4/11/2013		23581	7723	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.7 U	0.00E+00	5.54E-04	116.97
Pulse-off period April 11, 2013 to May 10, 2013																
5/10/2013		23583	7724	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.5 U	0.00E+00	9.56E-04	116.97
7/15/2013		25160	8039	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.6 U	0.00E+00	7.61E-04	117.21
7/15/2013	Dup	25160	8039	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.7 U	0.00E+00	7.43E-04	-
Pulse-off period July 15, 2013 to September 9, 2013																
9/9/2013		25162	8040	160	2.0 U	0.00E+00	2.0 U	0.00E+00	2.0 U	0.00E+00	20 U	0.00E+00	8.0 U	0.00E+00	1.86E-03	117.21
11/18/2013		26825	8372	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.6 U	0.00E+00	2.66E-04	117.30
Pulse-off period November 18, 2013 to January 15, 2014																
1/15/2014		28218	8651	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	7.44E-04	117.51
3/14/2014		29432	8894	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	5.75E-05	117.52
3/14/2014	Dup	29432	8894	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	5.0 U	0.00E+00	8.30E-05	-
Pulse-off period March 14, 2014 to May 15, 2014																
5/15/2014		29914	8990	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	25	3.60E-05	11 U	0.00E+00	1.21E-03	117.64
7/23/2014		31567	9321	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.6 U	0.00E+00	4.64E-04	117.79
Pulse-off period July 23, 2014 to September 16, 2014																
9/16/2014		32432	9494	160	2.1 U	0.00E+00	2.1 U	0.00E+00	2.1 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	1.51E-03	118.05
11/14/2014		33847	9777	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.6 U	0.00E+00	2.31E-04	118.12
Pulse-off period November 14, 2014 to January 9, 2015																
1/9/2015		33855	9778	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.8 U	0.00E+00	9.45E-04	118.12
3/13/2015		35189	10045	160	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	13 U	0.00E+00	5.0 U	0.00E+00	9.70E-05	118.15
Pulse-off period March 13, 2015 to May 15, 2015																
5/15/2015		35194	10046	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.6 U	0.00E+00	1.06E-03	118.15
7/16/2015		36677	10343	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.6 U	0.00E+00	3.62E-04	118.25
Pulse-off period July 16, 2015 to September 22, 2015																
9/22/2015		36680	10343	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.5 U	0.00E+00	2.25E-03	118.26

Notes:

Mass removal rate = (flow rate in scfm)(concentration in ppmv)(60)(MW) / (387*1000000)

"U" indicates non-detection at the specified reporting limit; for ND compounds, zero is used in mass removal calculations.

MW molecular weight (values from the U.S. National Library of Medicine)

SCFM standard cubic feet per minute

J Indicates estimated value.

B The analyte was detected in the method, field and/or trip blank.

When a duplicate sample was collected, the original sample results are used in the mass calculations.

Table 4.3
Cell 3 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 3 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 3 Run Time (hr)	SVE Flow Rate (scfm)	1,1,1-Trichloroethane		1,1,2-Trichloroethane		1,1-Dichloroethane		1,2-Dichloroethane		1,1-Dichloroethene		cis-1,2-Dichloroethene		trans-1,2-Dichloroethene		Tetrachloroethene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
12/14/2009		181	60	140	94000	2.72E-01	270 U	0.00E+00	1100	2.36E-03	270 U	0.00E+00	2300	4.84E-03	8100	1.70E-02	270 U	0.00E+00	750	2.70E-03
12/16/2009		229	76	150	46000	1.43E-01	110 U	0.00E+00	710	1.63E-03	110 U	0.00E+00	1100	2.48E-03	5500	1.24E-02	110 U	0.00E+00	400	1.54E-03
1/5/2010		707	236	140	42000	1.22E-01	150 U	0.00E+00	290	6.23E-04	150 U	0.00E+00	980	2.06E-03	1500	3.16E-03	150 U	0.00E+00	240	8.64E-04
1/21/2010		1084	361	150	15000	4.65E-02	42 U	0.00E+00	260	5.98E-04	42 U	0.00E+00	280	6.31E-04	1600	3.61E-03	42 U	0.00E+00	170	6.56E-04
1/21/2010	Dup	1084	361	150	16000	4.96E-02	43 U	0.00E+00	280	6.44E-04	43 U	0.00E+00	290	6.54E-04	1700	3.83E-03	43 U	0.00E+00	170	6.56E-04
2/24/2010		1893	631	150	11000	3.41E-02	28 U	0.00E+00	240	5.52E-04	28 U	0.00E+00	280	6.31E-04	1100	2.48E-03	28 U	0.00E+00	140	5.40E-04
3/15/2010		2345	782	140	20000	5.79E-02	21 U	0.00E+00	400	8.59E-04	21 U	0.00E+00	510	1.07E-03	1900	4.00E-03	21 U	0.00E+00	280	1.01E-03
4/14/2010		2804	935	150	31000	9.62E-02	100 U	0.00E+00	380	8.75E-04	100 U	0.00E+00	1100	2.48E-03	1200	2.71E-03	100 U	0.00E+00	820	3.16E-03
5/13/2010		3495	1165	140	8300	2.40E-02	12 U	0.00E+00	220	4.73E-04	12 U	0.00E+00	190	4.00E-04	960	2.02E-03	12 U	0.00E+00	200	7.20E-04
6/21/2010		4430	1477	108	7200	1.61E-02	21 U	0.00E+00	220	3.65E-04	21 U	0.00E+00	150	2.43E-04	660	1.07E-03	21 U	0.00E+00	160	4.44E-04
7/21/2010		5058	1686	140	6100	1.77E-02	20 U	0.00E+00	120	2.58E-04	20 U	0.00E+00	130	2.74E-04	460	9.68E-04	20 U	0.00E+00	120	4.32E-04
8/23/2010		5784	1928	0	8000	0.00E+00	20 U	0.00E+00	200	0.00E+00	20 U	0.00E+00	120	0.00E+00	490	0.00E+00	20 U	0.00E+00	220	0.00E+00
9/23/2010		6523	2174	145	6600	1.98E-02	11 U	0.00E+00	140	3.11E-04	11 U	0.00E+00	140	3.05E-04	440	9.59E-04	11 U	0.00E+00	160	5.96E-04
10/22/2010		7219	2406	140	3700	1.07E-02	14 U	0.00E+00	91	1.95E-04	14 U	0.00E+00	66	1.39E-04	210	4.42E-04	14 U	0.00E+00	110	3.96E-04
11/15/2010		7794	2598	140	4600	1.33E-02	15 U	0.00E+00	120	2.58E-04	15 U	0.00E+00	64	1.35E-04	170	3.58E-04	15 U	0.00E+00	88	3.17E-04
12/22/2010		8508	2777	150	5600	1.74E-02	20 U	0.00E+00	150	3.45E-04	20 U	0.00E+00	120	2.71E-04	330	7.44E-04	20 U	0.00E+00	56	2.16E-04
1/24/2011		9302	2975	170	2200	7.74E-03	8.3 U	0.00E+00	130	3.39E-04	8.3 U	0.00E+00	27	6.90E-05	200	5.11E-04	8.3 U	0.00E+00	35	1.53E-04
2/25/2011		10071	3167	165	1300	4.44E-03	4.0 U	0.00E+00	45	1.14E-04	4.0 U	0.00E+00	25	6.20E-05	72	1.79E-04	4.0 U	0.00E+00	28	1.19E-04
3/18/2011		10573	3293	165	360	1.23E-03	1.3 U	0.00E+00	24	6.08E-05	1.3 U	0.00E+00	5.4	1.34E-05	35	8.68E-05	1.3 U	0.00E+00	13	5.51E-05
4/15/2011		11241	3460	160	160 J,B	5.29E-04	1.0 U	0.00E+00	17	4.17E-05	1.0 U	0.00E+00	2.8	6.73E-06	28	6.73E-05	1.0 U	0.00E+00	15	6.17E-05
5/19/2011		12061	3665	160	64	2.12E-04	1.2 U	0.00E+00	10	2.45E-05	1.2 U	0.00E+00	1.4	3.37E-06	12	2.89E-05	1.2 U	0.00E+00	9.6	3.95E-05
6/16/2011		12722	3830	170	160	5.63E-04	1.2 U	0.00E+00	280	7.30E-04	1.2 U	0.00E+00	2.5	6.39E-06	34	8.69E-05	1.2 U	0.00E+00	61	2.67E-04
7/15/2011		13417	4472	170	190	6.68E-04	1.2 U	0.00E+00	8.3	2.16E-05	1.2 U	0.00E+00	2.8	7.15E-06	23	5.88E-05	1.2 U	0.00E+00	22	9.62E-05
8/22/2011		14324	4775	170	1600	5.63E-03	4.3 U	0.00E+00	19	4.96E-05	4.3 U	0.00E+00	21	5.37E-05	130	3.32E-04	4.3 U	0.00E+00	39	1.70E-04
9/15/2011		14905	4968	170	800	2.81E-03	3.7 U	0.00E+00	9.5	2.48E-05	3.7 U	0.00E+00	12	3.07E-05	62	1.58E-04	3.7 U	0.00E+00	24	1.05E-04
10/14/2011		15598	5199	160	750	2.48E-03	3.0 U	0.00E+00	10	2.45E-05	3.0 U	0.00E+00	13	3.13E-05	37	8.90E-05	3.0 U	0.00E+00	15	6.17E-05
11/21/2011		16510	5503	170	380	1.34E-03	1.4 U	0.00E+00	6.6	1.72E-05	1.4 U	0.00E+00	5.6	1.43E-05	24	6.13E-05	1.4 U	0.00E+00	7.9	3.45E-05
12/14/2011		17010	5670	170	830	2.92E-03	3.5 U	0.00E+00	8.7	2.27E-05	3.5 U	0.00E+00	70	1.79E-04	33	8.43E-05	3.5 U	0.00E+00	6.9	3.02E-05
1/19/2012		17923	5974	170	800	2.81E-03	3.0 U	0.00E+00	12	3.13E-05	3.0 U	0.00E+00	13	3.32E-05	33	8.43E-05	3.0 U	0.00E+00	6.4	2.80E-05
2/15/2012		18566	6189	170	1600	5.63E-03														

Table 4.3
Cell 3 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 3 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 3 Run Time (hr)	SVE Flow Rate (scfm)	Trichloroethene		Vinyl chloride		Methylene Chloride		Carbon Tetrachloride		Chloroform		Chloroethane		Benzene		Toluene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
12/14/2009		181	60	140	1000	2.85E-03	270 U	0.00E+00	270 U	0.00E+00	270 U	0.00E+00	270 U	0.00E+00	270 U	0.00E+00	270 U	0.00E+00	270 U	0.00E+00
12/16/2009		229	76	150	550	1.68E-03	110 U	0.00E+00	110 U	0.00E+00	110 U	0.00E+00	110 U	0.00E+00	110 U	0.00E+00	110 U	0.00E+00	110 U	0.00E+00
1/5/2010		707	236	140	250	7.13E-04	150 U	0.00E+00	220	4.06E-04	150 U	0.00E+00	150 U	0.00E+00	150 U	0.00E+00	150 U	0.00E+00	150 U	0.00E+00
1/21/2010		1084	361	150	140	4.28E-04	42 U	0.00E+00	42 U	0.00E+00	43 U	0.00E+00	43 U	0.00E+00	42 U	0.00E+00	42 U	0.00E+00	42 U	0.00E+00
1/21/2010	Dup	1084	361	150	140	4.28E-04	43 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00
2/24/2010		1893	631	150	66	2.02E-04	28 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00
3/15/2010		2345	782	140	120	3.42E-04	51	6.92E-05	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00
4/14/2010		2804	935	150	190	5.81E-04	100 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00
5/13/2010		3495	1165	140	43	1.23E-04	12 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00
6/21/2010		4430	1477	108	55	1.21E-04	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00
7/21/2010		5058	1686	140	44	1.25E-04	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00
8/23/2010		5784	1928	0	66	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00
9/23/2010		6523	2174	145	50	1.48E-04	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00
10/22/2010		7219	2406	140	31	8.84E-05	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00
11/15/2010		7794	2598	140	29	8.27E-05	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00
12/22/2010		8508	2777	150	21	6.42E-05	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00
1/24/2011		9302	2975	170	17	5.89E-05	8.3 U	0.00E+00	8.3 U	0.00E+00	8.3 U	0.00E+00	8.3 U	0.00E+00	8.3 U	0.00E+00	8.3 U	0.00E+00	8.3 U	0.00E+00
2/25/2011		10071	3167	165	16	5.38E-05	4.0 U	0.00E+00	4.0 U	0.00E+00	4.0 U	0.00E+00	4.0 U	0.00E+00	16 U	0.00E+00	4.0 U	0.00E+00	4.0 U	0.00E+00
3/18/2011		10573	3293	165	5.9	1.98E-05	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	5.4 U	0.00E+00	1.9	3.80E-06	1.3 U	0.00E+00
4/15/2011		11241	3460	160	7.7	2.51E-05	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	4.1 U	0.00E+00	2.6	5.04E-06	1.0 U	0.00E+00
5/19/2011		12061	3665	160	6.9	2.25E-05	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.8 U	0.00E+00	1.9	3.68E-06	1.2 U	0.00E+00
6/16/2011		12722	3830	170	9.8	3.39E-05	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.6	3.29E-06	1.2 U	0.00E+00
7/15/2011		13417	4472	170	9.3	3.22E-05	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.8 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
8/22/2011		14324	4775	170	21	7.27E-05	4.3 U	0.00E+00	4.3 U	0.00E+00	4.3 U	0.00E+00	4.3 U	0.00E+00	17 U	0.00E+00	4.3 U	0.00E+00	4.3 U	0.00E+00
9/15/2011		14905	4968	170	14	4.85E-05	3.7 U	0.00E+00	3.7 U	0.00E+00	3.7 U	0.00E+00	3.7 U	0.00E+00	15 U	0.00E+00	4.1	8.44E-06	3.7 U	0.00E+00
10/14/2011		15598	5199	160	13	4.24E-05	3.0 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00	12 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00
11/21/2011		16510	5503	170	9.2	3.19E-05	1.4 U	0.00E+00	1.4 UJ	0.00E+00	1.4 U	0.00E+00	1.4 U	0.00E+00	5.5 U	0.00E+00	1.4 U	0.00E+00	1.4 U	0.00E+00
12/14/2011		17010	5670	170	22	7.62E-05	3.5 U	0.00E+00	35 U	0.00E+00	3.5 U	0.00E+00	3.5 U	0.00E+00	14 U	0.00E+00	3.5 U	0.00E+00	3.5 U	0.00E+00
1/19/2012		17923	5974	170	12	4.16E-05	3.0 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00	12 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00
2/15/2012		18566	6189	170	24	8.31E-05	4.5 U	0.00E+00	4.5 U	0.00E+00	4.5 U</td									

Table 4.3
Cell 3 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 3 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 3 Run Time (hr)	SVE Flow Rate (scfm)	Ethylbenzene		m&p-Xylenes		o-Xylenes		Acetone		Methyl Ethyl Ketone (MEK)		Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)		
12/14/2009	Dup	181	60	140	270 U	0.00E+00	1600	3.69E-03	510	1.18E-03	1100 U	0.00E+00	270 U	0.00E+00	3.07E-01	18.51
12/16/2009		229	76	150	110 U	0.00E+00	540	1.33E-03	240	5.93E-04	590	7.97E-04	110 U	0.00E+00	1.65E-01	21.16
1/5/2010		707	236	140	150 U	0.00E+00	150 U	0.00E+00	150 U	0.00E+00	590 U	0.00E+00	150 U	0.00E+00	1.29E-01	41.78
1/21/2010		1084	361	150	42 U	0.00E+00	42 U	0.00E+00	42 U	0.00E+00	170 U	0.00E+00	42 U	0.00E+00	5.25E-02	48.37
1/21/2010		1084	361	150	43 U	0.00E+00	43 U	0.00E+00	43 U	0.00E+00	170 U	0.00E+00	43 U	0.00E+00	5.59E-02	48.80
2/24/2010		1893	631	150	28 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00	110 U	0.00E+00	28 U	0.00E+00	3.85E-02	58.76
3/15/2010		2345	782	140	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	83 U	0.00E+00	21 U	0.00E+00	6.53E-02	68.60
4/14/2010		2804	935	150	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	420 U	0.00E+00	100 U	0.00E+00	1.06E-01	84.81
5/13/2010		3495	1165	140	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	48 U	0.00E+00	12 U	0.00E+00	2.78E-02	91.21
6/21/2010		4430	1477	108	21 U	0.00E+00	21 U	0.00E+00	21 U	0.00E+00	83 U	0.00E+00	21 U	0.00E+00	1.83E-02	96.92
7/21/2010		5058	1686	140	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	79 U	0.00E+00	20 U	0.00E+00	1.97E-02	101.05
8/23/2010		5784	1928	0	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	81 U	0.00E+00	20 U	0.00E+00	0.00E+00	101.05
9/23/2010		6523	2174	145	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	43 U	0.00E+00	11 U	0.00E+00	2.21E-02	106.49
10/22/2010		7219	2406	140	14 U	0.00E+00	14 U	0.00E+00	14 U	0.00E+00	55 U	0.00E+00	14 U	0.00E+00	1.20E-02	109.27
11/15/2010		7794	2598	140	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	59 U	0.00E+00	15 U	0.00E+00	1.45E-02	112.05
12/22/2010		8508	2777	150	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	82 U	0.00E+00	20 U	0.00E+00	1.90E-02	115.44
1/24/2011		9302	2975	170	8.3 U	0.00E+00	8.3 U	0.00E+00	8.3 U	0.00E+00	33 U	0.00E+00	8.3 U	0.00E+00	8.87E-03	117.20
2/25/2011		10071	3167	165	4.0 U	0.00E+00	4.0 U	0.00E+00	4.0 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	4.96E-03	118.15
3/18/2011		10573	3293	165	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	10	1.49E-05	5.4 U	0.00E+00	1.48E-03	118.34
4/15/2011		11241	3460	160	1.0 U	0.00E+00	1.0 U	0.00E+00	1.0 U	0.00E+00	7.3 J,B	1.05E-05	4.1 U	0.00E+00	7.48E-04	118.47
5/19/2011		12061	3665	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.8 U	0.00E+00	4.8 U	0.00E+00	3.34E-04	118.53
6/16/2011		12722	3830	170	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	6.8	1.04E-05	4.7 U	0.00E+00	1.70E-03	118.81
7/15/2011		13417	4472	170	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	7.7	1.18E-05	4.8 U	0.00E+00	8.96E-04	119.39
8/22/2011		14324	4775	170	4.3 U	0.00E+00	4.3 U	0.00E+00	4.3 U	0.00E+00	17 U	0.00E+00	17 U	0.00E+00	6.30E-03	121.30
9/15/2011		14905	4968	170	3.7 U	0.00E+00	3.7 U	0.00E+00	3.7 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	3.19E-03	121.91
10/14/2011		15598	5199	160	3.0 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	2.73E-03	122.54
11/21/2011		16510	5503	170	1.4 U	0.00E+00	1.4 U	0.00E+00	1.4 U	0.00E+00	5.5 U,J	0.00E+00	5.5 U	0.00E+00	1.50E-03	123.00
12/14/2011		17010	5670	170	3.5 U	0.00E+00	3.5 U	0.00E+00	3.5 U	0.00E+00	380 J	5.82E-04	58	1.10E-04	4.00E-03	123.67
1/19/2012		17923	5974	170	3.0 U	0.00E+00	3.0 U	0.00E+00	3.0 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	3.03E-03	124.59
2/15/2012		18566	6189	170	4.5 U	0.00E+00	4.5 U	0.00E+00	4.5 U	0.00E+00	18 U	0.00E+00	18 U	0.00E+00	6.70E-03	126.03
3/15/2012		19262	6421	170	5.1 U	0.00E+00	5.1 U	0.00E+00	5.1 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	6.04E-03	127.43
4/19/2012		20102	6701	160	1.8 U	0.00E+00	1.8 U	0.00E+00	1.8 U	0.00E+00	7.3 U	0.00E+00	7.3 U	0.00E+00	2.13E-03	128.02
5/16/2012		20748	6916	160	0.80 U	0.00E+00	0.80 U	0.00E+00	0.80 U	0.00E+00	3.2 U	0.00E+00	3.2 U	0.00E+00	1.16E-03	128.27

See notes on last page.

2015 Q3 Tables 4.1 - 4.6 Air Tbl 4.3 C-3

Table 4.3
Cell 3 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 3 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 3 Run Time (hr)	SVE Flow Rate (scfm)	1,1,1-Trichloroethane		1,1,2-Trichloroethane		1,1-Dichloroethane		1,2-Dichloroethane		1,1-Dichloroethene		cis-1,2-Dichloroethene		trans-1,2-Dichloroethene		Tetrachloroethene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
Pulse-off period June 1, 2012 to August 14, 2012																				
8/14/2012		21282	7094	160	1200	3.97E-03	4.7 U	0.00E+00	20	4.91E-05	4.7 U	0.00E+00	17	4.09E-05	8.4	2.02E-05	4.7 U	0.00E+00	26	1.07E-04
9/17/2012		21952	7317	160	5000	1.65E-02	16 U	0.00E+00	38	9.33E-05	16 U	0.00E+00	16 U	0.00E+00	130	3.13E-04	16 U	0.00E+00	100	4.11E-04
Pulse-off period September 17, 2012 to November 15, 2012																				
11/15/2012		21959	7320	160	640	2.12E-03	2.1 U	0.00E+00	130	3.19E-04	2.1 U	0.00E+00	22	5.29E-05	18	4.33E-05	2.1 U	0.00E+00	18	7.40E-05
12/14/2012		22554	7518	170	1100	3.87E-03	4.6 U	0.00E+00	17 J	4.43E-05	4.6 U	0.00E+00	15	3.83E-05	70 J	1.79E-04	4.6 U	0.00E+00	7.1 J	3.10E-05
12/14/2012	Dup	22554	7518	170	1100 J	3.87E-03	4.5 U	0.00E+00	25 J	6.52E-05	4.5 U	0.00E+00	17 J	4.34E-05	36 J	9.20E-05	4.5 U	0.00E+00	49 J	2.14E-04
Pulse-off period December 14, 2012 to February 26, 2013																				
2/26/2013		22556	7519	160	640	2.12E-03	2.4 U	0.00E+00	12	2.95E-05	2.4 U	0.00E+00	8	1.92E-05	23	5.53E-05	2.4 U	0.00E+00	15	6.17E-05
4/11/2013		23581	8134	160	180	5.96E-04	1.2 U	0.00E+00	7	1.72E-05	1.2 U	0.00E+00	2.8	6.73E-06	28	6.73E-05	1.2 U	0.00E+00	4.6	1.89E-05
Pulse-off period April 11, 2013 to May 10, 2013																				
5/10/2013		23583	8135	160	200	6.62E-04	1.1 U	0.00E+00	5.5	1.35E-05	1.1 U	0.00E+00	3.3	7.94E-06	22	5.29E-05	1.1 U	0.00E+00	4.4	1.81E-05
7/15/2013		25160	9082	160	64	2.12E-04	1.2 U	0.00E+00	3.6	8.84E-06	1.2 U	0.00E+00	1.2 U	0.00E+00	1.9	4.57E-06	1.2 U	0.00E+00	14	5.76E-05
Pulse-off period July 15, 2013 to September 9, 2013																				
9/9/2013		25162	9083	160	390	1.29E-03	1.7 U	0.00E+00	5.6	1.37E-05	1.7 U	0.00E+00	3	9.93E-06	2	4.81E-06	1.7 U	0.00E+00	21	8.64E-05
11/18/2013		26825	10081	160	350	1.16E-03	1.2 U	0.00E+00	8.2	2.01E-05	1.2 U	0.00E+00	6.6	2.18E-05	45	1.08E-04	1.2 U	0.00E+00	5.7	2.34E-05
Pulse-off period November 18, 2013 to January 15, 2014																				
1/15/2014		28218	10916	160	240	7.94E-04	1.2 U	0.00E+00	5	1.23E-05	1.2 U	0.00E+00	4.1	1.36E-05	16	3.85E-05	1.2 U	0.00E+00	18	7.40E-05
3/14/2014		29432	11645	160	72	2.38E-04	1.2 U	0.00E+00	8.7	2.14E-05	1.2 U	0.00E+00	2.4	7.94E-06	6.4	1.54E-05	1.2 U	0.00E+00	9.5	3.91E-05
Pulse-off period March 14, 2014 to May 15, 2014																				
5/15/2014		29914	11934	160	770	2.55E-03	2.3 U	0.00E+00	15	3.68E-05	2.3 U	0.00E+00	12	3.97E-05	86	2.07E-04	2.3 U	0.00E+00	6.9	2.84E-05
7/23/2014		31567	12926	160	130	4.30E-04	1.4 U	0.00E+00	5	1.23E-05	1.4 U	0.00E+00	1.4	4.63E-06	10	2.40E-05	1.4 U	0.00E+00	10	4.11E-05
Pulse-off period July 23, 2014 to September 16, 2014																				
9/16/2014		32432	13099	160	390	1.29E-03	2.4 U	0.00E+00	15	3.68E-05	2.4 U	0.00E+00	3	7.21E-06	8.4	2.02E-05	2.4 U	0.00E+00	17	6.99E-05
11/14/2014		33847	13948	160	180	5.96E-04	1.2 U	0.00E+00	5.2	1.28E-05	1.2 U	0.00E+00	3	9.93E-06	25	6.01E-05	1.2 U	0.00E+00	18	7.40E-05
Pulse-off period November 14, 2014 to January 9, 2015																				
1/9/2015		33855	13949	160	220	7.28E-04	1.1 U	0.00E+00	4.7	1.15E-05	1.1 U	0.00E+00	2.2	5.29E-06	18	4.33E-05	1.1 U	0.00E+00	11	4.52E-05
3/13/2015		35189	14750	160	200	6.62E-04	1.2 U	0.00E+00	4.4	1.08E-05	1.2 U	0.00E+00	3.1	1.03E-05	14	3.37E-05	1.2 U	0.00E+00	4.2	1.73E-05
Pulse-off period March 13, 2015 to May 15, 2015																				
5/15/2015		35194	14751	160	300	9.93E-04	1.2 U	0.00E+00	5.6	1.37E-05	1.2 U	0.00E+00	3.1	7.45E-06	10	2.40E-05	1.2 U	0.00E+00	8.1	3.33E-05
7/16/2015		36677	15641	160	180	5.96E-04	1.2 U	0.00E+00	6.5	1.60										

Table 4.3
Cell 3 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 3 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 3 Run Time (hr)	SVE Flow Rate (scfm)	Trichloroethene		Vinyl chloride		Methylene Chloride		Carbon Tetrachloride		Chloroform		Chloroethane		Benzene		Toluene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
Pulse-off period June 1, 2012 to August 14, 2012																				
8/14/2012		21282	7094	160	25	8.15E-05	4.7 U	0.00E+00	47 U	0.00E+00	4.7 U	0.00E+00	4.7 U	0.00E+00	19 U	0.00E+00	4.7 U	0.00E+00	4.7 U	0.00E+00
9/17/2012		21952	7317	160	180	5.87E-04	16 U	0.00E+00	160 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	65 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00
Pulse-off period September 17, 2012 to November 15, 2012																				
11/15/2012		21959	7320	160	39	1.27E-04	2.1 U	0.00E+00	21 U	0.00E+00	2.1 U	0.00E+00	2.1 U	0.00E+00	8.5 U	0.00E+00	2.1 U	0.00E+00	2.1 U	0.00E+00
12/14/2012		22554	7518	170	90 J	3.12E-04	4.6 U	0.00E+00	46 U	0.00E+00	4.6 U	0.00E+00	4.6 U	0.00E+00	18 U	0.00E+00	4.6 U	0.00E+00	4.6 U	0.00E+00
12/14/2012	Dup	22554	7518	170	52 J	1.80E-04	4.5 U	0.00E+00	45 U	0.00E+00	4.5 U	0.00E+00	4.5 U	0.00E+00	18 U	0.00E+00	4.5 U	0.00E+00	4.5 U	0.00E+00
Pulse-off period December 14, 2012 to February 26, 2013																				
2/26/2013		22556	7519	160	36	1.17E-04	2.4 U	0.00E+00	24 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	9.5 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00
4/11/2013		23581	8134	160	15	4.89E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
Pulse-off period April 11, 2013 to May 10, 2013																				
5/10/2013		23583	8135	160	13	4.24E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.3 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00
7/15/2013		25160	9082	160	10	3.26E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
Pulse-off period July 15, 2013 to September 9, 2013																				
9/9/2013		25162	9083	160	12	3.91E-05	1.7 U	0.00E+00	17 U	0.00E+00	1.7 U	0.00E+00	1.7 U	0.00E+00	6.8 U	0.00E+00	1.7 U	0.00E+00	1.7 U	0.00E+00
11/18/2013		26825	10081	160	14	4.56E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
Pulse-off period November 18, 2013 to January 15, 2014																				
1/15/2014		28218	10916	160	7.6	2.48E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.6 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
3/14/2014		29432	11645	160	8.1	2.64E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.9 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
Pulse-off period March 14, 2014 to May 15, 2014																				
5/15/2014		29914	11934	160	20	6.52E-05	2.3 U	0.00E+00	23 U	0.00E+00	2.3 U	0.00E+00	2.3 U	0.00E+00	9.3 U	0.00E+00	2.3 U	0.00E+00	2.3 U	0.00E+00
7/23/2014		31567	12926	160	9	2.93E-05	1.4 U	0.00E+00	14 U	0.00E+00	1.4 U	0.00E+00	1.4 U	0.00E+00	5.6 U	0.00E+00	1.4 U	0.00E+00	1.4 U	0.00E+00
Pulse-off period July 23, 2014 to September 16, 2014																				
9/16/2014		32432	13099	160	14	4.56E-05	2.4 U	0.00E+00	24 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	9.5 U	0.00E+00	3	5.81E-06	2.4 U	0.00E+00
11/14/2014		33847	13948	160	6.2	2.02E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.6 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
Pulse-off period November 14, 2014 to January 9, 2015																				
1/9/2015		33855	13949	160	6	1.96E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.3 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00
3/13/2015		35189	14750	160	14	4.56E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.6 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
Pulse-off period March 13, 2015 to May 15, 2015																				
5/15/2015		35194	14751	160	10	3.26E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
7/16/2015</																				

Table 4.3
Cell 3 - Phase 1 SVE System Effluent Data
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 3 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 3 Run Time (hr)	SVE Flow Rate (scfm)	Ethylbenzene		m&p-Xylenes		o-Xylenes		Acetone		Methyl Ethyl Ketone (MEK)		Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)		
Pulse-off period June 1, 2012 to August 14, 2012																
8/14/2012		21282	7094	160	4.7 U	0.00E+00	4.7 U	0.00E+00	4.7 U	0.00E+00	47 U	0.00E+00	19 U	0.00E+00	4.27E-03	129.03
9/17/2012		21952	7317	160	16 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	160 U	0.00E+00	65 U	0.00E+00	1.80E-02	133.04
Pulse-off period September 17, 2012 to November 15, 2012																
11/15/2012		21959	7320	160	2.1 U	0.00E+00	2.1 U	0.00E+00	2.1 U	0.00E+00	21 U	0.00E+00	8.5 U	0.00E+00	2.73E-03	133.05
12/14/2012		22554	7518	170	4.6 U	0.00E+00	4.6 U	0.00E+00	4.6 U	0.00E+00	46 U	0.00E+00	18 U	0.00E+00	4.47E-03	133.94
12/14/2012	Dup	22554	7518	170	4.5 U	0.00E+00	4.5 U	0.00E+00	4.5 U	0.00E+00	45 U	0.00E+00	18 U	0.00E+00	4.46E-03	-
Pulse-off period December 14, 2012 to February 26, 2013																
2/26/2013		22556	7519	160	2.4 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	24 U	0.00E+00	9.5 U	0.00E+00	2.40E-03	133.94
4/11/2013		23581	8134	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.7 U	0.00E+00	7.55E-04	134.40
Pulse-off period April 11, 2013 to May 10, 2013																
5/10/2013		23583	8135	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.3 U	0.00E+00	7.97E-04	134.40
7/15/2013		25160	9082	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.7 U	0.00E+00	3.15E-04	134.70
Pulse-off period July 15, 2013 to September 9, 2013																
9/9/2013		25162	9083	160	1.7 U	0.00E+00	1.7 U	0.00E+00	1.7 U	0.00E+00	17 U	0.00E+00	6.8 U	0.00E+00	1.44E-03	134.70
11/18/2013		26825	10081	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.7 U	0.00E+00	1.38E-03	136.08
Pulse-off period November 18, 2013 to January 15, 2014																
1/15/2014		28218	10916	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.6 U	0.00E+00	9.57E-04	136.88
3/14/2014		29432	11645	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.9 U	0.00E+00	3.48E-04	137.13
Pulse-off period March 14, 2014 to May 15, 2014																
5/15/2014		29914	11934	160	2.3 U	0.00E+00	2.3 U	0.00E+00	2.3 U	0.00E+00	23 U	0.00E+00	9.3 U	0.00E+00	2.92E-03	137.98
7/23/2014		31567	12926	160	1.4 U	0.00E+00	1.4 U	0.00E+00	1.4 U	0.00E+00	14 U	0.00E+00	5.6 U	0.00E+00	5.42E-04	138.52
Pulse-off period July 23, 2014 to September 16, 2014																
9/16/2014		32432	13099	160	2.4 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	24 U	0.00E+00	9.5 U	0.00E+00	1.48E-03	138.77
11/14/2014		33847	13948	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12	1.73E-05	4.6 U	0.00E+00	7.90E-04	139.44
Pulse-off period November 14, 2014 to January 9, 2015																
1/9/2015		33855	13949	160	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.3 U	0.00E+00	8.53E-04	139.44
3/13/2015		35189	14750	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.6 U	0.00E+00	7.79E-04	140.07
Pulse-off period March 13, 2015 to May 15, 2015																
5/15/2015		35194	14751	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.7 U	0.00E+00	1.10E-03	140.07
7/16/2015		36677	15641	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.9 U	0.00E+00	7.29E-04	140.72
Pulse-off period July 16, 2015 to September 22, 2015																
9/22/2015		36680	15641	160	2.3 U	0.00E+00	2.3 U	0.00E+00	2.3 U	0.00E+00	23 U	0.00E+00	9.3 U	0.00E+00	1.93E-03	140.72

Notes:

Mass removal rate = (flow rate in scfm)(concentration in ppmv)(60)(MW) / (387*1000000)

"U" indicates non-detection at the specified reporting limit; for ND compounds, zero is used in mass removal calculations.

MW = Molecular weight
SCFM = standard cubic feet per minute

J = Indicates estimated value.

B = The analyte was detected in the method, field and/or trip blank.

When a duplicate sample was collected, the original sample results are used in the mass calculations.

Table 4.4
Cell 4 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 4 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 4 Run Time (hr)	SVE Flow Rate (scfm)	1,1,1-Trichloroethane		1,1,2-Trichloroethane		1,1-Dichloroethane		1,2-Dichloroethane		1,1-Dichloroethene		cis-1,2-Dichloroethene		trans-1,2-Dichloroethene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
3/11/2011		222	222	500	150000	1.55E+00	600 U	0.00E+00	1800	1.38E-02	600 U	0.00E+00	860	6.46E-03	1400	1.05E-02	600 U	0.00E+00
3/18/2011		366	366	500	41000	4.24E-01	150 U	0.00E+00	1000	7.67E-03	150 U	0.00E+00	250	1.88E-03	460	3.46E-03	150 U	0.00E+00
3/18/2011	Dup	366	366	500	40000	4.14E-01	130 U	0.00E+00	1000	7.67E-03	130 U	0.00E+00	300	2.25E-03	480	3.61E-03	130 U	0.00E+00
3/25/2011		463	463	500	22000	2.28E-01	62 U	0.00E+00	980	7.52E-03	62 U	0.00E+00	87	6.54E-04	290	2.18E-03	62 U	0.00E+00
3/30/2011		558	558	500	25000	2.59E-01	68 U	0.00E+00	820	6.29E-03	68 U	0.00E+00	190	1.43E-03	290	2.18E-03	68 U	0.00E+00
4/8/2011		764	764	500	22000	2.28E-01	80 U	0.00E+00	1000	7.67E-03	80 U	0.00E+00	170	1.28E-03	340	2.56E-03	80 U	0.00E+00
4/15/2011		924	924	500	18000	1.86E-01	84 U	0.00E+00	930	7.13E-03	84 U	0.00E+00	110	8.27E-04	280	2.10E-03	84 U	0.00E+00
4/15/2011	Dup	924	924	500	16000 J	1.65E-01	60 U	0.00E+00	820 J	6.29E-03	60 U	0.00E+00	60 UJ	0.00E+00	260 J	1.95E-03	60 U	0.00E+00
5/19/2011		1685	1685	500	11000	1.14E-01	11 U	0.00E+00	640	4.91E-03	11 U	0.00E+00	100	7.52E-04	190	1.43E-03	11 U	0.00E+00
6/16/2011		2191	2191	420	10000	8.69E-02	11 U	0.00E+00	530	3.42E-03	11 U	0.00E+00	110 J	6.94E-04	160	1.01E-03	11 U	0.00E+00
6/16/2011	Dup	2191	2191	420	9600	8.34E-02	11 U	0.00E+00	510	3.29E-03	11 U	0.00E+00	110 J	6.94E-04	160	1.01E-03	11 U	0.00E+00
7/15/2011		2750	2750	420	7600	6.60E-02	24 U	0.00E+00	290	1.87E-03	24 U	0.00E+00	58	3.66E-04	79	4.99E-04	24 U	0.00E+00
8/22/2011		3133	3133	420	9000	7.82E-02	27 U	0.00E+00	410	2.64E-03	27 U	0.00E+00	92	5.81E-04	160	1.01E-03	27 U	0.00E+00
8/22/2011	Dup	3133	3133	420	9000	7.82E-02	22 U	0.00E+00	400	2.58E-03	22 U	0.00E+00	80	5.05E-04	150	9.47E-04	22 U	0.00E+00
9/15/2011		3630	3630	420	7000	6.08E-02	22 U	0.00E+00	250	1.61E-03	22 U	0.00E+00	55	3.47E-04	97	6.12E-04	22 U	0.00E+00
10/14/2011		4226	4226	420	4400	3.82E-02	19 U	0.00E+00	180	1.16E-03	19 U	0.00E+00	59	3.72E-04	60	3.79E-04	19 U	0.00E+00
11/21/2011		5019	5019	380	3700	2.91E-02	16 U	0.00E+00	170	9.91E-04	16 U	0.00E+00	320	1.83E-03	63	3.60E-04	16 U	0.00E+00
12/14/2011		5343	5343	260	4000	2.15E-02	19 U	0.00E+00	140	5.58E-04	19 U	0.00E+00	300	1.17E-03	55	2.15E-04	19 U	0.00E+00
1/19/2012		5993	5993	0	5200	0.00E+00	24 U	0.00E+00	160	0.00E+00	24 U	0.00E+00	58	0.00E+00	38	0.00E+00	24 U	0.00E+00
2/15/2012		6368	6368	260	4200	2.26E-02	19 U	0.00E+00	100	3.99E-04	19 U	0.00E+00	700	2.74E-03	53	2.07E-04	19 U	0.00E+00
3/15/2012		6946	6946	350	4000	2.90E-02	15 U	0.00E+00	120	6.44E-04	15 U	0.00E+00	38	2.00E-04	38	2.00E-04	15 U	0.00E+00
4/19/2012		7629	7629	380	5200	4.09E-02	16 U	0.00E+00	160	9.33E-04	16 U	0.00E+00	42	2.40E-04	43	2.46E-04	16 U	0.00E+00
5/16/2012		8143	8143	420	4100	3.56E-02	15 U	0.00E+00	110	7.09E-04	15 U	0.00E+00	43	2.71E-04	40	2.53E-04	15 U	0.00E+00

Table 4.4
Cell 4 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 4 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 4 Run Time (hr)	SVE Flow Rate (scfm)	Tetrachloroethene		Trichloroethene		Vinyl chloride		Methylene Chloride		Carbon Tetrachloride		Chloroform		Chloroethane		Benzene	
					Conc (ppbv)	Mass (lb/hr)	Conc (ppbv)	Mass (lb/hr)	Conc (ppbv)	Mass (lb/hr)	Conc (ppbv)	Mass (lb/hr)	Conc (ppbv)	Mass (lb/hr)	Conc (ppbv)	Mass (lb/hr)	Conc (ppbv)	Mass (lb/hr)	Conc (ppbv)	Mass (lb/hr)
3/11/2011		222	222	500	7200	9.26E-02	3900	3.97E-02	600 U	0.00E+00	600 U	0.00E+00	600 U	0.00E+00	600 U	0.00E+00	2400 U	0.00E+00	600 U	0.00E+00
3/18/2011		366	366	500	2900	3.73E-02	1600	1.63E-02	150 U	0.00E+00	150 U	0.00E+00	150 U	0.00E+00	150 U	0.00E+00	750 J	3.75E-03	150 U	0.00E+00
3/18/2011	Dup	366	366	500	3000	3.86E-02	1600	1.63E-02	130 UJ	0.00E+00	130 U	0.00E+00	130 U	0.00E+00	130 U	0.00E+00	1100 J	5.50E-03	130 U	0.00E+00
3/25/2011		463	463	500	3200	4.11E-02	970	9.88E-03	62 U	0.00E+00	61 NJ	4.02E-04	62 U	0.00E+00	62 U	0.00E+00	610	3.05E-03	62 U	0.00E+00
3/30/2011		558	558	500	2500	3.21E-02	1000	1.02E-02	68 U	0.00E+00	68 U	0.00E+00	68 U	0.00E+00	68 U	0.00E+00	470	2.35E-03	68 U	0.00E+00
4/8/2011		764	764	500	2400	3.09E-02	1000	1.02E-02	80 U	0.00E+00	80 U	0.00E+00	80 U	0.00E+00	80 U	0.00E+00	430	2.15E-03	80 U	0.00E+00
4/15/2011		924	924	500	1700	2.19E-02	920	9.37E-03	84 U	0.00E+00	84 U	0.00E+00	84 U	0.00E+00	84 U	0.00E+00	340 U	0.00E+00	84 U	0.00E+00
4/15/2011	Dup	924	924	500	1500 J	1.93E-02	830 J	8.45E-03	60 U	0.00E+00	60 U	0.00E+00	60 U	0.00E+00	60 U	0.00E+00	260 J	1.30E-03	60 U	0.00E+00
5/19/2011		1685	1685	500	1400	1.80E-02	530	5.40E-03	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	67	3.35E-04	26	1.57E-04
6/16/2011		2191	2191	420	1000	1.08E-02	410	3.51E-03	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	46 U	0.00E+00	14	7.12E-05
6/16/2011	Dup	2191	2191	420	960	1.04E-02	400	3.42E-03	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	45 U	0.00E+00	12	6.10E-05
7/15/2011		2750	2750	420	570	6.16E-03	250	2.14E-03	24 U	0.00E+00	28	1.55E-04	24 U	0.00E+00	24 U	0.00E+00	95 U	0.00E+00	24 U	0.00E+00
8/22/2011		3133	3133	420	920	9.93E-03	380	3.25E-03	27 U	0.00E+00	27 U	0.00E+00	27 U	0.00E+00	27 U	0.00E+00	110 U	0.00E+00	27 U	0.00E+00
8/22/2011	Dup	3133	3133	420	940	1.02E-02	360	3.08E-03	22 U	0.00E+00	22 U	0.00E+00	22 U	0.00E+00	22 U	0.00E+00	90 U	0.00E+00	22 U	0.00E+00
9/15/2011		3630	3630	420	660	7.13E-03	270	2.31E-03	22 U	0.00E+00	22 U	0.00E+00	22 U	0.00E+00	22 U	0.00E+00	90 U	0.00E+00	22 U	0.00E+00
10/14/2011		4226	4226	420	390	4.21E-03	180	1.54E-03	19 U	0.00E+00	19 U	0.00E+00	19 U	0.00E+00	19 U	0.00E+00	77 U	0.00E+00	19 U	0.00E+00
11/21/2011		5019	5019	380	360	3.52E-03	180	1.39E-03	16 U	0.00E+00	160 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	63 U	0.00E+00	16 U	0.00E+00
12/14/2011		5343	5343	260	360	2.41E-03	160	8.47E-04	19 U	0.00E+00	190 U	0.00E+00	19 U	0.00E+00	19 U	0.00E+00	74 U	0.00E+00	19 U	0.00E+00
1/19/2012		5993	5993	0	320	0.00E+00	180	0.00E+00	24 U	0.00E+00	24 U	0.00E+00	24 U	0.00E+00	24 U	0.00E+00	97 U	0.00E+00	24 U	0.00E+00
2/15/2012		6368	6368	260	280	1.87E-03	150	7.94E-04	19 U	0.00E+00	19 U	0.00E+00	19 U	0.00E+00	19 U	0.00E+00	78 U	0.00E+00	19 U	0.00E+00
3/15/2012		6946	6946	350	240	2.16E-03	140	9.98E-04	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	58 U	0.00E+00	15 U	0.00E+00
4/19/2012		7629	7629	380	400	3.91E-03	180	1.39E-03	16 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	62 U	0.00E+00	16 U	0.00E+00
5/16/2012		8143	8143	420	320	3.46E-03	150	1.28E-03	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	61 U	0.00E+00	15 U	0.00E+00

Table 4.4
Cell 4 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 4 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 4 Run Time (hr)	SVE Flow Rate (scfm)	Toluene		Ethylbenzene		m&p-Xylenes		o-Xylenes		Acetone		Methyl-Ethyl Ketone (MEK)		Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)		
3/11/2011		222	222	500	600 U	0.00E+00	600 U	0.00E+00	710	5.84E-03	600 U	0.00E+00	2400 U	0.00E+00	2400 U	0.00E+00	1.72E+00	381.87
3/18/2011		366	366	500	620 J	4.43E-03	150 U	0.00E+00	240	1.98E-03	200	1.65E-03	1500 J	6.75E-03	590 U	0.00E+00	5.09E-01	453.50
3/18/2011	Dup	366	366	500	380 J	2.71E-03	130 U	0.00E+00	250	2.06E-03	240	1.98E-03	690 J	3.11E-03	540 U	0.00E+00	4.97E-01	453.50
3/25/2011		463	463	500	140	1.00E-03	62 U	0.00E+00	78	6.42E-04	67	5.51E-04	250 U	0.00E+00	250 U	0.00E+00	2.95E-01	482.07
3/30/2011		558	558	500	190	1.36E-03	68 U	0.00E+00	250	2.06E-03	140	1.15E-03	270 U	0.00E+00	270 U	0.00E+00	3.18E-01	512.25
4/8/2011		764	764	500	200	1.43E-03	120	9.88E-04	560	4.61E-03	260	2.14E-03	320 U	0.00E+00	320 U	0.00E+00	2.91E-01	572.27
4/15/2011		924	924	500	170	1.21E-03	110	9.05E-04	540	4.44E-03	260	2.14E-03	340 U	0.00E+00	340 U	0.00E+00	2.36E-01	610.05
4/15/2011	Dup	924	924	500	140 J	1.00E-03	99 J	8.15E-04	540 J	4.44E-03	230 J	1.89E-03	240 J,B	1.08E-03	240 U	0.00E+00	2.12E-01	610.05
5/19/2011		1685	1685	500	100	7.14E-04	140	1.15E-03	920	7.57E-03	420	3.46E-03	81	3.65E-04	43 U	0.00E+00	1.58E-01	730.28
6/16/2011		2191	2191	420	51	3.06E-04	83	5.74E-04	600	4.15E-03	280	1.94E-03	46 J,B	1.74E-04	46 U	0.00E+00	1.14E-01	753.86
6/16/2011	Dup	2191	2191	420	53	3.18E-04	78	5.39E-04	580	4.01E-03	270	1.87E-03	69 J,B	2.61E-04	45 U	0.00E+00	1.09E-01	785.55
7/15/2011		2750	2750	420	28	1.68E-04	41	2.83E-04	270	1.87E-03	120	8.30E-04	180	6.81E-04	95 U	0.00E+00	8.10E-02	830.85
8/22/2011		3133	3133	420	35 J	2.10E-04	59 J	4.08E-04	340	2.35E-03	140	9.68E-04	110 U	0.00E+00	110 U	0.00E+00	9.95E-02	868.97
8/22/2011	Dup	3133	3133	420	22 UJ	0.00E+00	30 J	2.07E-04	310	2.14E-03	130	8.99E-04	90 U	0.00E+00	90 U	0.00E+00	9.87E-02	868.65
9/15/2011		3630	3630	420	22 U	0.00E+00	31	2.14E-04	340	2.35E-03	130	8.99E-04	90 U	0.00E+00	90 U	0.00E+00	7.63E-02	906.88
10/14/2011		4226	4226	420	38	2.28E-04	19 U	0.00E+00	170	1.18E-03	70	4.84E-04	77 U	0.00E+00	77 U	0.00E+00	4.78E-02	935.35
11/21/2011		5019	5019	380	16 U	0.00E+00	17	1.06E-04	220	1.38E-03	100	6.25E-04	160 U	0.00E+00	63 U	0.00E+00	3.93E-02	966.50
12/14/2011		5343	5343	260	19 U	0.00E+00	19 U	0.00E+00	76	3.25E-04	55	2.35E-04	190 UJ	0.00E+00	74 U	0.00E+00	2.73E-02	975.34
1/19/2012		5993	5993	0	36	0.00E+00	24 U	0.00E+00	78	0.00E+00	50	0.00E+00	97 U	0.00E+00	97 U	0.00E+00	0.00E+00	975.34
2/15/2012		6368	6368	260	19 U	0.00E+00	19 U	0.00E+00	58	2.48E-04	40	1.71E-04	300	7.02E-04	78 U	0.00E+00	2.97E-02	986.48
3/15/2012		6946	6946	350	15 U	0.00E+00	15 U	0.00E+00	44	2.53E-04	31	1.79E-04	58 U	0.00E+00	58 U	0.00E+00	3.36E-02	1005.89
4/19/2012		7629	7629	380	16 U	0.00E+00	16 U	0.00E+00	48	3.00E-04	33	2.06E-04	62 U	0.00E+00	62 U	0.00E+00	4.81E-02	1038.74
5/16/2012		8143	8143	420	15 U	0.00E+00	15 U	0.00E+00	28	1.94E-04	23	1.59E-04	61 U	0.00E+00	61 U	0.00E+00	4.19E-02	1060.30

Table 4.4
Cell 4 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 4 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 4 Run Time (hr)	SVE Flow Rate (scfm)	1,1,1-Trichloroethane		1,1,2-Trichloroethane		1,1-Dichloroethane		1,2-Dichloroethane		1,1-Dichloroethene		cis-1,2-Dichloroethene		trans-1,2-Dichloroethene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
Pulse-off period June 1, 2012 to August 14, 2012																		
8/14/2012		8546	8546	420	5000	4.34E-02	16 U	0.00E+00	98	6.32E-04	16 U	0.00E+00	66	4.17E-04	27	1.70E-04	16 U	0.00E+00
9/17/2012		9033	9033	470	3700	3.60E-02	15 U	0.00E+00	140	1.01E-03	15 U	0.00E+00	15 U	0.00E+00	26	1.84E-04	15 U	0.00E+00
Pulse-off period September 17, 2012 to November 15, 2012																		
11/15/2012		9037	9037	420	4900 J	4.26E-02	28 U	0.00E+00	74 J	4.77E-04	28 U	0.00E+00	110 J	6.94E-04	29 J	1.83E-04	28 U	0.00E+00
11/15/2012	Dup	9037	9037	420	8700	7.56E-02	24 U	0.00E+00	200 J	1.29E-03	24 U	0.00E+00	220	1.39E-03	360 J	2.27E-03	24 U	0.00E+00
12/14/2012		9439	9439	150	500	1.55E-03	1.9 U	0.00E+00	14	3.22E-05	1.9 U	0.00E+00	6.8	1.53E-05	18	4.06E-05	1.9 U	0.00E+00
Pulse-off period December 14, 2012 to February 26, 2013																		
2/26/2013		9439	9439	0	520	0.00E+00	2.2 U	0.00E+00	23	0.00E+00	2.2 U	0.00E+00	5.7	0.00E+00	28	0.00E+00	2.2 U	0.00E+00
4/11/2013		9876	9876	340	430	3.02E-03	1.8 U	0.00E+00	26	1.36E-04	1.8 U	0.00E+00	7.1	3.63E-05	28	1.43E-04	1.8 U	0.00E+00
Pulse-off period April 11, 2013 to May 10, 2013																		
5/10/2013		9882	9882	340	270	1.90E-03	1.1 U	0.00E+00	23	1.20E-04	1.1 U	0.00E+00	3.4	1.74E-05	30	1.53E-04	1.1 U	0.00E+00
7/15/2013		10907	10907	340	100	7.03E-04	1.1 U	0.00E+00	13	6.78E-05	1.1 U	0.00E+00	1.7	8.69E-06	14	7.15E-05	1.1 U	0.00E+00
Pulse-off period July 15, 2013 to September 9, 2013																		
9/9/2013		10914	10914	340	170	1.20E-03	1.2 U	0.00E+00	17	8.87E-05	1.2 U	0.00E+00	2.2	1.12E-05	27	1.38E-04	1.2	6.13E-06
11/18/2013		11992	11992	260	330	1.77E-03	1.1 U	0.00E+00	7.9	3.15E-05	1.1 U	0.00E+00	5.2	2.03E-05	14	5.47E-05	1.1 U	0.00E+00
Pulse-off period November 18, 2013 to March 14, 2014																		
1/15/2014		11997	11997	320	200	1.32E-03	1.2 U	0.00E+00	5.5	2.70E-05	1.2 U	0.00E+00	3.3	1.59E-05	9.6	4.62E-05	1.2 U	0.00E+00
3/14/2014		12980	12980	180	430	1.60E-03	2.6 U	0.00E+00	6.2	1.71E-05	2.6 U	0.00E+00	8.2	2.22E-05	18	4.87E-05	2.6 U	0.00E+00
Pulse-off period March 14, 2014 to May 15, 2014																		
5/15/2014		12986	12986	180	470	1.75E-03	1.1 U	0.00E+00	10	2.76E-05	1.1 U	0.00E+00	6.9	1.87E-05	22	5.95E-05	1.1 U	0.00E+00
7/23/2014		14627	14627	300	14	8.69E-05	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	1.6	0.00E+00	1.3 U	0.00E+00
Pulse-off period July 23, 2014 to September 16, 2014																		
9/16/2014		14634	14628	320	150	9.93E-04	1.2 U	0.00E+00	9	4.42E-05	1.2 U	0.00E+00	1.7	8.18E-06	15	7.21E-05	1.2 U	0.00E+00
11/14/2014		16008	16008	320	220	1.46E-03	0.96 U	0.00E+00	5	2.45E-05	0.96 U	0.00E+00	3.6	1.73E-05	8.9	4.28E-05	0.96 U	0.00E+00
Pulse-off period November 14, 2014 to January 9, 2015																		
1/9/2015		16015	16015	260	150	8.07E-04	1.1 U	0.00E+00	4.1	1.64E-05	1.1 U	0.00E+00	2.2	8.60E-06	7.4	2.89E-05	1.1 U	0.00E+00
3/13/2015		17178	17178	220	190	8.65E-04	1.2 U	0.00E+00	4.9	1.65E-05	1.2 U	0.00E+00	3.1	1.03E-05	5.5	1.82E-05	1.2 U	0.00E+00
Pulse-off period March 13, 2015 to May 15, 2015																		
5/15/2015		17186	17186	320	180	1.19E-03	2.6 U	0.00E+00	4.3	2.11E-05	2.6 U	0.00E+00	2.8	1.35E-05	5.2	2.50E-05	2.6 U	0.00E+00
7/16/2015		18436	18436	310	270	1.73E-03	1.2 U	0.00E+00	7.7	3.66E-05	1.2 U	0.00E+00	4	1.86E-05	13	6.06E-05	1.2 U	0.00E+00
Pulse-off period July 16, 2015 to September 22, 2015																		
9/22/2015		18439	18439	300	200	1.24E-03	1.1 U	0.00E+00	6.3	2.90E-05	1.1 U	0.00E+00	2.1	9.47E-06	11	4.96E-05	1.1 U	0.00E+00

Notes:

Mass removal rate = (flow rate in scfm)(concentration in ppmv)(60)(MW) / (387*1000000)

Table 4.4
Cell 4 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 4 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 4 Run Time (hr)	SVE Flow Rate (scfm)	Tetrachloroethene		Trichloroethene		Vinyl chloride		Methylene Chloride		Carbon Tetrachloride		Chloroform		Chloroethane		Benzene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
Pulse-off period June 1, 2012 to August 14, 2012																				
8/14/2012		8546	8546	420	490	5.29E-03	180	1.54E-03	16 U	0.00E+00	160 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	63 U	0.00E+00	16 U	0.00E+00
9/17/2012		9033	9033	470	410	4.95E-03	220	2.11E-03	15 U	0.00E+00	150 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	61 U	0.00E+00	15 U	0.00E+00
Pulse-off period September 17, 2012 to November 15, 2012																				
11/15/2012	Dup	9037	9037	420	260 J	2.81E-03	150 J	1.28E-03	28 U	0.00E+00	280 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00	110 U	0.00E+00	28 U	0.00E+00
11/15/2012		9037	9037	420	1200 J	1.30E-02	390 J	3.34E-03	24 U	0.00E+00	240 U	0.00E+00	24 U	0.00E+00	24 U	0.00E+00	94 U	0.00E+00	24 U	0.00E+00
12/14/2012		9439	9439	150	62	2.39E-04	28	8.56E-05	1.9 U	0.00E+00	19 U	0.00E+00	1.9 U	0.00E+00	1.9 U	0.00E+00	7.5 U	0.00E+00	1.9 U	0.00E+00
Pulse-off period December 14, 2012 to February 26, 2013																				
2/26/2013		9439	9439	0	130	0.00E+00	27	0.00E+00	2.2 U	0.00E+00	22 U	0.00E+00	2.2 U	0.00E+00	2.2 U	0.00E+00	8.7 U	0.00E+00	2.2 U	0.00E+00
4/11/2013		9876	9876	340	98	8.57E-04	25	1.73E-04	1.8 U	0.00E+00	18 U	0.00E+00	1.8 U	0.00E+00	1.8 U	0.00E+00	7.1 U	0.00E+00	1.8 U	0.00E+00
Pulse-off period April 11, 2013 to May 10, 2013																				
5/10/2013		9882	9882	340	120	1.05E-03	23	1.59E-04	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.5 U	0.00E+00	1.1 U	0.00E+00
7/15/2013		10907	10907	340	180	1.57E-03	30	2.08E-04	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.6 U	0.00E+00	1.1 U	0.00E+00
Pulse-off period July 15, 2013 to September 9, 2013																				
9/9/2013		10914	10914	340	350	3.06E-03	50	3.46E-04	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.6 U	0.00E+00	1.2 U	0.00E+00
11/18/2013		11992	11992	260	50	3.34E-04	13	6.89E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.4 U	0.00E+00	1.1 U	0.00E+00
Pulse-off period November 18, 2013 to March 14, 2014																				
1/15/2014		11997	11997	320	51	4.20E-04	11	7.17E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.8 U	0.00E+00	1.2 U	0.00E+00
3/14/2014		12980	12980	180	7.8	3.61E-05	14	5.13E-05	2.6 U	0.00E+00	26 U	0.00E+00	2.6 U	0.00E+00	2.6 U	0.00E+00	10 U	0.00E+00	2.6 U	0.00E+00
Pulse-off period March 14, 2014 to May 15, 2014																				
5/15/2014		12986	12986	180	38	1.76E-04	17	6.23E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.6 U	0.00E+00	1.1 U	0.00E+00
7/23/2014		14627	14627	300	15	1.16E-04	2.4	1.47E-05	1.3 U	0.00E+00	13 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	5.2 U	0.00E+00	1.3 U	0.00E+00
Pulse-off period July 23, 2014 to September 16, 2014																				
9/16/2014		14634	14628	320	200	1.65E-03	39	2.54E-04	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.9 U	0.00E+00	2	7.75E-06
11/14/2014		16008	16008	320	69	5.68E-04	12	7.82E-05	0.96 U	0.00E+00	9.6 U	0.00E+00	0.96 U	0.00E+00	0.96 U	0.00E+00	3.8 U	0.00E+00	0.96 U	0.00E+00
Pulse-off period November 14, 2014 to January 9, 2015																				
1/9/2015		16015	16015	260	50	3.34E-04	11	5.83E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.4 U	0.00E+00	1.1 U	0.00E+00
3/13/2015		17178	17178	220	27	1.53E-04	6.9	3.09E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.8 U	0.00E+00	1.2 U	0.00E+00
Pulse-off period March 13, 2015 to May 15, 2015																				
5/15/2015		17186	17186	320	45	3.70E-04	9.8	6.39E-05	2.6 U	0.00E+00	26 U	0.00E+00	2.6 U	0.00E+00	2.6 U	0.00E+00	10 U	0.00E+00	2.6 U	0.00E+00
7/16/2015		18436	18436	310	130	1.04E-03	27	1.71E-04	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U							

Table 4.4
Cell 4 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 4 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 4 Run Time (hr)	SVE Flow Rate (scfm)	Toluene		Ethylbenzene		m&p-Xylenes		o-Xylenes		Acetone		Methyl Ethyl Ketone (MEK)		Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)		
Pulse -off period	June 1, 2012 to August 14, 2012																	
8/14/2012		8546	8546	420	16 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	160 U	0.00E+00	63 U	0.00E+00	5.15E-02	1081.05
9/17/2012		9033	9033	470	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	150 U	0.00E+00	61 U	0.00E+00	4.42E-02	1102.58
Pulse -off period	September 17, 2012 to November 15, 2012																	
11/15/2012		9037	9037	420	28 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00	28 U	0.00E+00	280 U	0.00E+00	110 U	0.00E+00	4.80E-02	1102.78
11/15/2012	Dup	9037	9037	420	24 U	0.00E+00	24 U	0.00E+00	24 U	0.00E+00	24 U	0.00E+00	240 U	0.00E+00	94 U	0.00E+00	9.68E-02	1102.97
12/14/2012		9439	9439	150	1.9 U	0.00E+00	1.9 U	0.00E+00	1.9 U	0.00E+00	1.9 U	0.00E+00	19 U	0.00E+00	7.5 U	0.00E+00	1.96E-03	1103.57
Pulse -off period	December 14, 2012 to February 26, 2013																	
2/26/2013		9439	9439	0	2.2 U	0.00E+00	2.2 U	0.00E+00	2.2 U	0.00E+00	2.2 U	0.00E+00	22 U	0.00E+00	8.7 U	0.00E+00	0.00E+00	1103.57
4/11/2013		9876	9876	340	1.8 U	0.00E+00	1.8 U	0.00E+00	1.8 U	0.00E+00	1.8 U	0.00E+00	18 U	0.00E+00	7.1 U	0.00E+00	4.37E-03	1105.48
Pulse -off period	April 11, 2013 to May 10, 2013																	
5/10/2013		9882	9882	340	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.5 U	0.00E+00	3.40E-03	1105.50
7/15/2013		10907	10907	340	19	9.23E-05	1.2	6.72E-06	2.2	1.23E-05	1.1 U	0.00E+00	24	7.35E-05	4.9	1.86E-05	2.84E-03	1108.40
Pulse -off period	July 15, 2013 to September 9, 2013																	
9/9/2013		10914	10914	340	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	14	4.29E-05	4.6 U	0.00E+00	4.89E-03	1108.44
11/18/2013		11992	11992	260	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.4 U	0.00E+00	2.28E-03	1110.90
Pulse -off period	November 18, 2013 to March 14, 2014																	
1/15/2014		11997	11997	320	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.8 U	0.00E+00	1.90E-03	1110.91
3/14/2014		12980	12980	180	2.6 U	0.00E+00	2.6 U	0.00E+00	2.6 U	0.00E+00	2.6 U	0.00E+00	26 U	0.00E+00	10 U	0.00E+00	1.78E-03	1112.65
Pulse -off period	March 14, 2014 to May 15, 2014																	
5/15/2014		12986	12986	180	3.9	1.00E-05	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.6 U	0.00E+00	2.10E-03	1112.67
7/23/2014		14627	14627	300	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	1.3 U	0.00E+00	13 U	0.00E+00	5.2 U	0.00E+00	2.17E-04	1113.02
Pulse -off period	July 23, 2014 to September 16, 2014																	
9/16/2014		14634	14628	320	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	21	6.05E-05	4.9 U	0.00E+00	3.09E-03	1113.03
11/14/2014		16008	16008	320	0.96 U	0.00E+00	0.96 U	0.00E+00	0.96 U	0.00E+00	0.96 U	0.00E+00	9.6 U	0.00E+00	3.8 U	0.00E+00	2.19E-03	1116.04
Pulse -off period	November 14, 2014 to January 9, 2015																	
1/9/2015		16015	16015	260	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.4 U	0.00E+00	1.25E-03	1116.05
3/13/2015		17178	17178	220	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.8 U	0.00E+00	1.09E-03	1117.32
Pulse -off period	March 13, 2015 to May 15, 2015																	
5/15/2015		17186	17186	320	2.6 U	0.00E+00	2.6 U	0.00E+00	2.6 U	0.00E+00	2.6 U	0.00E+00	26 U	0.00E+00	10 U	0.00E+00	1.68E-03	1117.34
7/16/2015		18436	18436	310	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.9 U	0.00E+00	3.05E-03	1121.16
Pulse -off period	July 16, 2015 to September 22, 2015																	
9/22/2015		18439	18439	300	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.3 U	0.00E+00	3.09E-03	1121.16

Notes:

Mass removal rate = (flow rate in scfm)(concentration in ppmv)(60)(MW) / (387*1000000)

"U" indicates non-detection at the specified reporting limit; for ND compounds, zero is used in mass removal calculations.

MW molecular weight (values from the U.S. National Library
SCFM standard cubic feet per minute

J Indicates estimated value.

B The analyte was detected in the method, field and/or trip blank.

When a duplicate sample was collected, the original sample results are used in the mass calculations.

Table 4.5
Cell 5 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 5 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 5 Run Time (hr)	SVE Flow Rate (scfm)	1,1,1-Trichloroethane		1,1,2-Trichloroethane		1,1-Dichloroethane		1,2-Dichloroethane		1,1-Dichloroethene		cis-1,2-Dichloroethene		trans-1,2-Dichloroethene		Tetrachloroethene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
3/11/2011		218	218	360	28000	2.08E-01	100 U	0.00E+00	2400	1.33E-02	100 U	0.00E+00	740	4.00E-03	10000	5.41E-02	100 U	0.00E+00	5900	5.46E-02
3/18/2011		362	362	360	13000	9.68E-02	52 U	0.00E+00	1100	6.08E-03	52 U	0.00E+00	280	1.52E-03	4800	2.60E-02	52 U	0.00E+00	6800	6.29E-02
3/25/2011		459	459	360	8900	6.63E-02	30 U	0.00E+00	650	3.59E-03	30 U	0.00E+00	200	1.08E-03	2600	1.41E-02	30 U	0.00E+00	5400	5.00E-02
3/30/2011		553	553	360	4600	3.43E-02	13 U	0.00E+00	310	1.71E-03	13 U	0.00E+00	100	5.41E-04	1300	7.03E-03	13 U	0.00E+00	4000	3.70E-02
4/8/2011		759	759	360	4600	3.43E-02	20 U	0.00E+00	330	1.82E-03	20 U	0.00E+00	95	5.14E-04	1100	5.95E-03	20 U	0.00E+00	5700	5.28E-02
4/15/2011		920	920	360	4600	3.43E-02	20 U	0.00E+00	370	2.04E-03	20 U	0.00E+00	69	3.73E-04	980	5.30E-03	20 U	0.00E+00	4600	4.26E-02
5/19/2011		1681	1681	330	2800	1.91E-02	12 U	0.00E+00	250	1.27E-03	12 U	0.00E+00	34	1.69E-04	730	3.62E-03	12 U	0.00E+00	7800	6.62E-02
6/16/2011		2187	2187	300	1800	1.12E-02	7.8 U	0.00E+00	170	7.82E-04	7.8 U	0.00E+00	23 J	1.04E-04	520	2.34E-03	7.8 U	0.00E+00	2400	1.85E-02
7/15/2011		2745	2745	220	2400	1.09E-02	7.6 U	0.00E+00	180	6.08E-04	7.6 U	0.00E+00	27	8.93E-05	840	2.78E-03	7.6 U	0.00E+00	2700	1.53E-02
8/22/2011		3129	3129	260	1700	9.14E-03	5.0 U	0.00E+00	150	5.98E-04	5.0 U	0.00E+00	21	8.21E-05	690	2.70E-03	5.0 U	0.00E+00	2000	1.34E-02
9/15/2011		3626	3626	220	1400	6.37E-03	4.5 U	0.00E+00	69	2.33E-04	4.5 U	0.00E+00	22	7.27E-05	380	1.26E-03	4.5 U	0.00E+00	1100	6.22E-03
10/14/2011		4222	4222	220	980	4.46E-03	3.9 U	0.00E+00	57	1.92E-04	3.9 U	0.00E+00	19	6.28E-05	310	1.03E-03	3.9 U	0.00E+00	760	4.30E-03
11/21/2011	Dup	5015	5015	200	690	2.85E-03	3.2 U	0.00E+00	55	1.69E-04	3.2 U	0.00E+00	45	1.35E-04	290	8.72E-04	3.2 U	0.00E+00	380	1.95E-03
11/21/2011	Dup	5015	5015	200	700	2.90E-03	3.1 U	0.00E+00	57	1.75E-04	3.1 U	0.00E+00	59	1.77E-04	300	9.02E-04	3.1 U	0.00E+00	390	2.01E-03
12/14/2011		5339	5339	200	890	3.68E-03	3.2 U	0.00E+00	62	1.90E-04	3.2 U	0.00E+00	64	1.92E-04	270	8.12E-04	3.2 U	0.00E+00	350	1.80E-03
1/19/2012		5958	5958	0	540	0.00E+00	2.8 U	0.00E+00	17	0.00E+00	2.8 U	0.00E+00	9.9	0.00E+00	69	0.00E+00	2.8 U	0.00E+00	78	0.00E+00
2/15/2012		6364	6364	0	990	0.00E+00	4.1 U	0.00E+00	24	0.00E+00	4.1 U	0.00E+00	100	0.00E+00	230	0.00E+00	4.1 U	0.00E+00	150	0.00E+00
3/15/2012		6942	6942	0	1100	0.00E+00	3.8 U	0.00E+00	43	0.00E+00	3.8 U	0.00E+00	20	0.00E+00	220	0.00E+00	3.8 U	0.00E+00	140	0.00E+00
4/19/2012		7625	7625	80	650	1.08E-03	2.4 U	0.00E+00	28	3.44E-05	2.4 U	0.00E+00	8.1	9.74E-06	130	1.56E-04	2.4 U	0.00E+00	100	2.06E-04
5/16/2012		8138	8138	200	650	2.69E-03	2.0 U	0.00E+00	28	8.59E-05	2.0 U	0.00E+00	8.9	2.68E-05	110	3.31E-04	2.0 U	0.00E+00	130	6.68E-04
Pulse-off period June 1, 2012 to August 14, 2012																				
8/14/2012		8541	8541	360	710	3.23E-03	2.5 U	0.00E+00	44	1.49E-04	2.5 U	0.00E+00	11	3.64E-05	110	3.64E-04	2.5 U	0.00E+00	540	3.05E-03
9/17/2012		9029	9029	360	2000	8.27E-03	8.0 U	0.00E+00	29	8.90E-05	8.0 U	0.00E+00	19	5.71E-05	42	1.26E-04	8.0 U	0.00E+00	190	9.77E-04
Pulse-off period September 17, 2012 to November 15, 2012																				
11/15/2012		9033	9033	220	1200	5.46E-03	4.4 U	0.00E+00	19	6.41E-05	4.4 U	0.00E+00	33	1.09E-04	8	2.65E-05	4.4 U	0.00E+00	55	3.11E-04
12/14/2012		9436	9436	200	1200	4.96E-03	4.8 U	0.00E+00	35	1.07E-04	4.8 U	0.00E+00	16	4.81E-05	37	1.11E-04	4.8 U	0.00E+00	61	3.14E-04
Pulse-off period December 14, 2012 to February 26, 2013																				
2/26/2013		9511	9511	440	70	6.37E-04	6.8 U	0.00E+00	6.8 U	0.00E+00	6.8 U	0.00E+00	6.8 U	0.00E+00	6.8 U	0.00E+00	6.8 U	0.00E+00	6.8 U	0.00E+00
4/11/2013		9952	9952	420	1600	1.39E-02	8	6.95E-05	160	1.03E-03	5.1 U	0.00E+00	20	1.26E-04	88	5.56E-04	5.1 U	0.00E+00	320	3.46E-03
Pulse-off period April 11, 2013 to May 10, 2013																				
5/10/2013		9958	9958	420	1200	1.04E-02	5.4 U	0.00E+00	86	5.54E-04	5.4 U	0.00E+00	12	7.57E-05	45	2.84E-04				

Table 4.5
Cell 5 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 5 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 5 Run Time (hr)	SVE Flow Rate (scfm)	Trichloroethene		Vinyl chloride		Methylene Chloride		Carbon Tetrachloride		Chloroform		Chloroethane		Benzene		Toluene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
3/11/2011		218	218	360	1400	1.03E-02	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	420 U	0.00E+00	100 U	0.00E+00	350	1.80E-03
3/18/2011		362	362	360	1100	8.07E-03	52 U	0.00E+00	52 U	0.00E+00	52 U	0.00E+00	52 U	0.00E+00	210 U	0.00E+00	52 U	0.00E+00	120 JB	6.17E-04
3/25/2011		459	459	360	760	5.57E-03	30 U	0.00E+00	33	1.56E-04	30 U	0.00E+00	30 U	0.00E+00	120 U	0.00E+00	30 U	0.00E+00	73	3.75E-04
3/30/2011		553	553	360	420	3.08E-03	13 U	0.00E+00	13 U	0.00E+00	13 U	0.00E+00	13 U	0.00E+00	51 U	0.00E+00	13 U	0.00E+00	37	1.90E-04
4/8/2011		759	759	360	560	4.11E-03	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	81 U	0.00E+00	20 U	0.00E+00	57	2.93E-04
4/15/2011		920	920	360	560	4.11E-03	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	20 U	0.00E+00	81 U	0.00E+00	20 U	0.00E+00	85	4.37E-04
5/19/2011		1681	1681	330	360	2.42E-03	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	47 U	0.00E+00	12 U	0.00E+00	120	5.66E-04
6/16/2011		2187	2187	300	180	1.10E-03	7.8 U	0.00E+00	7.8 U	0.00E+00	7.8 U	0.00E+00	7.8 U	0.00E+00	31 U	0.00E+00	12	4.36E-05	7.8 U	0.00E+00
7/15/2011		2745	2745	220	280	1.25E-03	7.6 U	0.00E+00	20	5.79E-05	7.6 U	0.00E+00	7.6 U	0.00E+00	30 U	0.00E+00	7.6 U	0.00E+00	49	1.54E-04
8/22/2011		3129	3129	260	160	8.47E-04	5.0 U	0.00E+00	5.0 U	0.00E+00	5.0 U	0.00E+00	5.0 U	0.00E+00	20 U	0.00E+00	7.6	2.39E-05	5.0 U	0.00E+00
9/15/2011		3626	3626	220	83	3.72E-04	4.5 U	0.00E+00	4.5 U	0.00E+00	4.5 U	0.00E+00	4.5 U	0.00E+00	18 U	0.00E+00	5	1.33E-05	4.5 U	0.00E+00
10/14/2011		4222	4222	220	50	2.24E-04	3.9 U	0.00E+00	3.9 U	0.00E+00	3.9 U	0.00E+00	3.9 U	0.00E+00	16 U	0.00E+00	3.9 U	0.00E+00	3.9 U	0.00E+00
11/21/2011	Dup	5015	5015	200	27	1.10E-04	3.2 U	0.00E+00	32 U	0.00E+00	3.2 U	0.00E+00	3.2 U	0.00E+00	13 U	0.00E+00	3.2 U	0.00E+00	3.2 U	0.00E+00
11/21/2011		5015	5015	200	28	1.14E-04	3.1 U	0.00E+00	31 U	0.00E+00	3.1 U	0.00E+00	3.1 U	0.00E+00	12 U	0.00E+00	3.1 U	0.00E+00	3.1 U	0.00E+00
12/14/2011		5339	5339	200	24	9.78E-05	3.2 U	0.00E+00	32 U	0.00E+00	3.2 U	0.00E+00	3.2 U	0.00E+00	13 U	0.00E+00	3.2 U	0.00E+00	3.2 U	0.00E+00
1/19/2012		5958	5958	0	10	0.00E+00	2.8 U	0.00E+00	2.8 U	0.00E+00	2.8 U	0.00E+00	2.8 U	0.00E+00	11 U	0.00E+00	2.8 U	0.00E+00	2.8 U	0.00E+00
2/15/2012		6364	6364	0	19	0.00E+00	4.1 U	0.00E+00	4.1 U	0.00E+00	4.1 U	0.00E+00	4.1 U	0.00E+00	16 U	0.00E+00	4.1 U	0.00E+00	4.1 U	0.00E+00
3/15/2012		6942	6942	0	25	0.00E+00	3.8 U	0.00E+00	3.8 U	0.00E+00	3.8 U	0.00E+00	3.8 U	0.00E+00	15 U	0.00E+00	3.8 U	0.00E+00	3.8 U	0.00E+00
4/19/2012		7625	7625	80	19	3.10E-05	2.4 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	9.4 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00
5/16/2012		8138	8138	200	24	9.78E-05	2.0 U	0.00E+00	2.0 U	0.00E+00	2.0 U	0.00E+00	2.0 U	0.00E+00	7.9 U	0.00E+00	2.0 U	0.00E+00	2.0 U	0.00E+00
Pulse -off period June 1, 2012 to August 14, 2012																				
8/14/2012		8541	8541	360	64	2.87E-04	2.5 U	0.00E+00	25 U	0.00E+00	2.5 U	0.00E+00	2.5 U	0.00E+00	9.9 U	0.00E+00	2.5 U	0.00E+00	2.5 U	0.00E+00
9/17/2012		9029	9029	360	71	2.89E-04	8.0 U	0.00E+00	80 U	0.00E+00	8.0 U	0.00E+00	8.0 U	0.00E+00	32 U	0.00E+00	8.0 U	0.00E+00	8.0 U	0.00E+00
Pulse -off period September 17, 2012 to November 15, 2012																				
11/15/2012		9033	9033	220	39	1.75E-04	4.4 U	0.00E+00	44 U	0.00E+00	4.4 U	0.00E+00	4.4 U	0.00E+00	18 U	0.00E+00	4.4 U	0.00E+00	4.4 U	0.00E+00
12/14/2012		9436	9436	200	60	2.44E-04	4.8 U	0.00E+00	48 U	0.00E+00	4.8 U	0.00E+00	4.8 U	0.00E+00	19 U	0.00E+00	4.8 U	0.00E+00	4.8 U	0.00E+00
Pulse -off period December 14, 2012 to February 26, 2013																				
2/26/2013		9511	9511	440	6.8 U	0.00E+00	6.8 U	0.00E+00	68 U	0.00E+00	6.8 U	0.00E+00	6.8 U	0.00E+00	27 U	0.00E+00	12	6.39E-05	6.8 U	0.00E+00
4/11/2013		9952	9952	420	110	9.41E-04	5.1 U	0.00E+00	51 U	0.00E+00	5.1 U	0.00E+00	5.1 U	0.00E+00	20 U	0.00E+00	5.1 U	0.00E+00	5.1 U	0.00E+00
Pulse -off period April 11, 2013 to May 10, 2013																				
5/10/2013		9958	9958	420	79	6.76E-04	5.4 U	0.00E+00	54 U	0.00E+00	5.4 U	0.00E+00	5.4 U	0.00E+00	22 U	0.00E+00	5.4 U	0.00E+00	5.4 U	0.0

Table 4.5
Cell 5 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 5 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 5 Run Time (hr)	SVE Flow Rate (scfm)	Ethylbenzene		m&p-Xylenes		o-Xylenes		Acetone		Methyl/Ethyl Ketone (MEK)		Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)			
3/11/2011	Dup	218	218	360	100 U	0.00E+00	100 U	0.00E+00	100 U	0.00E+00	420 U	0.00E+00	420 U	0.00E+00	3.47E-01	75.54	
3/18/2011		362	362	360	52 U	0.00E+00	59	3.50E-04	110	6.52E-04	210 U	0.00E+00	210 U	0.00E+00	2.03E-01	104.77	
3/25/2011		459	459	360	30 U	0.00E+00	30 U	0.00E+00	47	2.79E-04	130	4.21E-04	120 U	0.00E+00	1.42E-01	118.53	
3/30/2011		553	553	360	16	9.48E-05	23	1.36E-04	46	2.73E-04	99	3.21E-04	51 U	0.00E+00	8.47E-02	126.48	
4/8/2011		759	759	360	38	2.25E-04	84	4.98E-04	120	7.11E-04	81 U	0.00E+00	81 U	0.00E+00	1.01E-01	147.32	
4/15/2011		920	920	360	45	2.67E-04	160	9.48E-04	140	8.30E-04	180 J,B	5.83E-04	81 U	0.00E+00	9.17E-02	162.08	
5/19/2011		1681	1681	330	12 U	0.00E+00	12 U	0.00E+00	12 U	0.00E+00	360	1.07E-03	47 U	0.00E+00	9.44E-02	233.92	
6/16/2011		2187	2187	300	15	7.41E-05	54	2.67E-04	64	3.16E-04	69 J,B	1.86E-04	31 U	0.00E+00	3.49E-02	251.58	
7/15/2011		2745	2745	220	13	4.71E-05	120	4.35E-04	140	5.07E-04	94	1.86E-04	30 U	0.00E+00	3.23E-02	269.61	
8/22/2011		3129	3129	260	5.9	2.52E-05	19	8.13E-05	29	1.24E-04	62 J,B	1.45E-04	20 U	0.00E+00	2.71E-02	280.03	
9/15/2011		3626	3626	220	4.5 U	0.00E+00	14	5.07E-05	17	6.16E-05	49	9.71E-05	18 U	0.00E+00	1.47E-02	287.36	
10/14/2011		4222	4222	220	3.9 U	0.00E+00	7.1	2.57E-05	10	3.62E-05	16 U	0.00E+00	16 U	0.00E+00	1.03E-02	293.51	
11/21/2011		5015	5015	200	3.2 U	0.00E+00	4.5	1.48E-05	6.1	2.01E-05	36	6.48E-05	13 U	0.00E+00	6.19E-03	298.43	
11/21/2011		5015	5015	200	3.1 U	0.00E+00	4.2	1.38E-05	6.2	2.04E-05	31 U	0.00E+00	12 U	0.00E+00	6.30E-03	298.51	
12/14/2011		5339	5339	200	3.2 U	0.00E+00	3.2 U	0.00E+00	3.2 U	0.00E+00	32 UJ	0.00E+00	13 U	0.00E+00	6.77E-03	300.62	
1/19/2012		5958	5958	0	2.8 U	0.00E+00	2.8 U	0.00E+00	2.8 U	0.00E+00	11 U	0.00E+00	11 U	0.00E+00	0.00E+00	300.62	
2/15/2012		6364	6364	0	4.1 U	0.00E+00	4.1 U	0.00E+00	4.1 U	0.00E+00	16 U	0.00E+00	16 U	0.00E+00	0.00E+00	300.62	
3/15/2012		6942	6942	0	3.8 U	0.00E+00	3.8 U	0.00E+00	3.8 U	0.00E+00	15 U	0.00E+00	15 U	0.00E+00	0.00E+00	300.62	
4/19/2012		7625	7625	80	2.4 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	9.4 U	0.00E+00	9.4 U	0.00E+00	1.51E-03	301.65	
5/16/2012		8138	8138	200	2.0 U	0.00E+00	2.0 U	0.00E+00	2.0 U	0.00E+00	7.9 U	0.00E+00	7.9 U	0.00E+00	3.90E-03	303.65	
Pulse-off period June 1, 2012 to August 14, 2012																	
8/14/2012		8541	8541	360	2.5 U	0.00E+00	2.5 U	0.00E+00	2.5 U	0.00E+00	25 U	0.00E+00	9.9 U	0.00E+00	7.12E-03	306.52	
9/17/2012		9029	9029	360	8.0 U	0.00E+00	8.0 U	0.00E+00	8.0 U	0.00E+00	80 U	0.00E+00	32 U	0.00E+00	9.81E-03	311.31	
Pulse-off period September 17, 2012 to November 15, 2012																	
11/15/2012		9033	9033	220	4.4 U	0.00E+00	4.4 U	0.00E+00	4.4 U	0.00E+00	44 U	0.00E+00	18 U	0.00E+00	6.15E-03	311.34	
12/14/2012		9436	9436	200	4.8 U	0.00E+00	4.8 U	0.00E+00	4.8 U	0.00E+00	48 U	0.00E+00	19 U	0.00E+00	5.79E-03	313.67	
Pulse-off period December 14, 2012 to February 26, 2013																	
2/26/2013		9511	9511	440	6.8 U	0.00E+00	6.8 U	0.00E+00	6.8 U	0.00E+00	68 U	0.00E+00	27 U	0.00E+00	7.01E-04	313.72	
4/11/2013		9952	9952	420	5.1 U	0.00E+00	5.1 U	0.00E+00	5.1 U	0.00E+00	51 U	0.00E+00	20 U	0.00E+00	2.01E-02	322.58	
Pulse-off period April 11, 2013 to May 10, 2013																	
5/10/2013		9958	9958	420	5.4 U	0.00E+00	5.4 U	0.00E+00	5.4 U	0.00E+00	54 U	0.00E+00	22 U	0.00E+00	1.44E-02	322.66	
7/15/2013		10984	10984	360	4.7 U	0.00E+00	4.7 U	0.00E+00	4.7 U	0.00E+00	47 U	0.00E+00	19 U	0.00E+00	1.65E-02	339.59	
Pulse-off period July 15, 2013 to September 9, 2013																	
9/9/2013		10991	10991	380	4 U	0.00E+00	4 U	0.00E+00	4 U	0.00E+00	40 U	0.00E+00	40 U	0.00E+00	8.81E-03	339.65	
11/18/2013		12069	12069	380	7.6 U	0.00E+00	7.6 U	0.00E+00	7.6 U	0.00E+00	76 U	0.00E+00	31 U	0.00E+00	1.58E-02	356.69	
Pulse-off period November 18, 2013 to January 15, 2014																	
1/15/2014		12074	12074	380	3.5 U	0.00E+00	3.5 U	0.00E+00	3.5 U	0.00E+00	35 U	0.00E+00	14 U	0.00E+00	8.88E-03	356.73	
3/14/2014		13057	13057	380	7.8 U	0.00E+00	7.8 U	0.00E+00	7.8 U	0.00E+00	78 U	0.00E+00	31 U	0.00E+00	1.24E-02	368.96	
Pulse-off period March 14, 2014 to May 15, 																	

Table 4.5
Cell 5 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 5 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 5 Run Time (hr)	SVE Flow Rate (scfm)	1,1,1-Trichloroethane		1,1,2-Trichloroethane		1,1-Dichloroethane		1,2-Dichloroethane		1,1-Dichloroethene		cis-1,2-Dichloroethene		trans-1,2-Dichloroethene		Tetrachloroethene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
Pulse-off period July 23, 2014 to September 16, 2014																				
9/16/2014		14721	14715	120	470	1.17E-03	2.3 U	0.00E+00	10	1.84E-05	2.3 U	0.00E+00	4.8	8.66E-06	6.9	1.24E-05	2.3 U	0.00E+00	79	2.44E-04
11/14/2014		16095	16095	290	660	3.96E-03	2.4 U	0.00E+00	15	6.67E-05	2.4 U	0.00E+00	8.5	3.70E-05	19	8.28E-05	2.4 U	0.00E+00	32	2.39E-04
Pulse-off period November 14, 2014 to January 9, 2015																				
1/9/2015		16102	16102	180	360	1.34E-03	1.1 U	0.00E+00	4.6	1.27E-05	1.1 U	0.00E+00	4.0	1.08E-05	7.2	1.95E-05	1.1 U	0.00E+00	12	5.55E-05
3/13/2015		17322	17322	260	660	3.55E-03	2.4 U	0.00E+00	22	8.78E-05	2.4 U	0.00E+00	8.0	3.13E-05	16	6.25E-05	2.4 U	0.00E+00	29	1.94E-04
Pulse-off period March 13, 2015 to May 15, 2015																				
5/15/2015		17329	17329	260	360	1.94E-03	1.1 U	0.00E+00	7.3	2.91E-05	1.1 U	0.00E+00	2.5	9.77E-06	5.9	2.31E-05	1.1 U	0.00E+00	31	2.07E-04
7/16/2015		18578	18578	180	260	9.68E-04	1.2 U	0.00E+00	22	6.08E-05	1.2 U	0.00E+00	3.5	9.47E-06	12	3.25E-05	1.2 U	0.00E+00	54	2.50E-04
Pulse-off period July 16, 2015 to September 22, 2015																				
9/22/2015		18580	18580	160	150	4.96E-04	1.2 U	0.00E+00	4.2	1.03E-05	1.2 U	0.00E+00	1.2	2.89E-06	2.4	5.77E-06	1.2 U	0.00E+00	47	1.93E-04

Notes:

Mass removal rate = (flow rate in scfm)(concentration in ppmv)(60)(MW) / (387*1000000)

"U" indicates non-detection at the specified reporting limit; for ND compounds, zero is used in mass removal calculations.

MW molecular weight (values from the U.S. National Library of Medicine)

SCFM standard cubic feet per minute

J Indicates estimated value.

B The analyte was detected in the method, field and/or trip blank.

When a duplicate sample was collected, the original sample results are used in the mass calculations.

Table 4.5
Cell 5 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 5 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 5 Run Time (hr)	SVE Flow Rate (scfm)	Trichloroethene		Vinyl chloride		Methylene Chloride		Carbon Tetrachloride		Chloroform		Chloroethane		Benzene		Toluene	
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)
Pulse-off period July 23, 2014 to September 16, 2014																				
9/16/2014		14721	14715	120	22	5.38E-05	2.3 U	0.00E+00	23 U	0.00E+00	2.3 U	0.00E+00	2.3 U	0.00E+00	9.4 U	0.00E+00	6.4	9.30E-06	2.3 U	0.00E+00
11/14/2014		16095	16095	290	11	6.50E-05	2.4 U	0.00E+00	24 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	9.7 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00
Pulse-off period November 14, 2014 to January 9, 2015																				
1/9/2015		16102	16102	180	4.9	1.80E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.6 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00
3/13/2015		17322	17322	260	12	6.36E-05	2.4 U	0.00E+00	24 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	9.5 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00
Pulse-off period March 13, 2015 to May 15, 2015																				
5/15/2015		17329	17329	260	8.2	4.34E-05	1.1 U	0.00E+00	11 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	4.5 U	0.00E+00	1.1 U	0.00E+00	1.4	5.20E-06
7/16/2015		18578	18578	180	14	5.13E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.7 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00
Pulse-off period July 16, 2015 to September 22, 2015																				
9/22/2015		18580	18580	160	11	3.59E-05	1.2 U	0.00E+00	12 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	4.9 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00

Notes:

Mass removal rate = (flow rate in scfm)(concentration in ppbv)(60)(MW) / (3871000000)

"U" indicates non-detection at the specified reporting limit; for ND compounds, zero is used in mass removal calculations.

MW molecular weight (values from the U.S. National Library of Medicine)

SCFM standard cubic feet per minute

J Indicates estimated value.

B The analyte was detected in the method, field and/or trip blank.

When a duplicate sample was collected, the original sample results are used in the mass calculations.

Table 4.5
Cell 5 - Phase 2 SVE System Effluent Data
March 2011 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

CELL 5 SVE EFFLUENT

Date	Sample Type	SVE Run Time (hr)	Cell 5 Run Time (hr)	SVE Flow Rate (scfm)	Ethylbenzene		m&p-Xylenes		o-Xylenes		Acetone		Methyl Ethyl Ketone (MEK)		Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)
					Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)	Conc (ppbv)	Mass Removal Rate (lb/hr)		
Pulse -off period July 23, 2014 to September 16, 2014																
9/16/2014		14721	14715	120	2.3 U	0.00E+00	2.3 U	0.00E+00	2.3 U	0.00E+00	32	3.46E-05	9.4 U	0.00E+00	1.55E-03	371.61
11/14/2014		16095	16095	290	2.4 U	0.00E+00	2.4 U	0.00E+00	2.8	1.34E-05	24 U	0.00E+00	9.7 U	0.00E+00	4.46E-03	377.77
Pulse -off period November 14, 2014 to January 9, 2015																
1/9/2015		16102	16102	180	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.6 U	0.00E+00	1.46E-03	377.78
3/13/2015		17322	17322	260	2.4 U	0.00E+00	2.4 U	0.00E+00	2.4 U	0.00E+00	24 U	0.00E+00	9.5 U	0.00E+00	3.99E-03	382.64
Pulse -off period March 13, 2015 to May 15, 2015																
5/15/2015		17329	17329	260	1.1 U	0.00E+00	1.1 U	0.00E+00	1.1 U	0.00E+00	11 U	0.00E+00	4.5 U	0.00E+00	2.25E-03	382.66
7/16/2015		18578	18578	180	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.7 U	0.00E+00	1.37E-03	384.37
Pulse -off period July 16, 2015 to September 22, 2015																
9/22/2015		18580	18580	160	1.2 U	0.00E+00	1.2 U	0.00E+00	1.2 U	0.00E+00	12 U	0.00E+00	4.9 U	0.00E+00	7.45E-04	384.37

Notes:

Mass removal rate = (flow rate in scfm)(concentration in ppmv)(60)(MW) / (387*1000000)

"U" indicates non-detection at the specified reporting limit; for ND compounds, zero is used in mass removal calculations.

MW molecular weight (values from the U.S. National Library of Medicine)

SCFM standard cubic feet per minute

J Indicates estimated value.

B The analyte was detected in the method, field and/or trip blank.

When a duplicate sample was collected, the original sample results are used in the mass calculations.

Table 4.6
Mass Removal - Phase 1 and Phase 2 AS/SVE Systems
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

Date	Cell 1			Cell 2			Cell 3			Cell 4			Cell 5			Total Cumulative Mass Removal (lb)
	Total Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	
12/3/2009	0															0.00
12/10/2009	53	0.22	11.91													11.91
12/11/2009				59	0.25	15.05										26.97
12/14/2009							60	0.31	18.51							45.48
12/15/2009				68	0.16	16.48										46.91
12/16/2009							76	0.17	21.16							49.55
12/22/2009	124	0.05	15.23													52.86
12/29/2009				180	0.12	29.76										66.15
1/5/2010							236	0.13	41.78							86.77
1/13/2010				301	0.05	35.75										92.75
1/21/2010							361	0.05	48.37							99.35
1/27/2010				408	0.06	42.68										106.27
2/24/2010	631	0.01	20.06	631	0.04	51.44	631	0.04	58.76							130.26
3/15/2010	782	0.01	22.02	782	0.09	64.40	782	0.07	68.60							155.02
4/14/2010	935	0.02	25.22	935	0.04	70.89	935	0.11	84.81							180.92
5/13/2010	1165	0.01	27.75	1165	0.04	79.74	1165	0.03	91.21							198.69
6/21/2010	1477	0.01	30.20	1477	0.02	86.90	1477	0.02	96.92							214.02
7/21/2010	1686	0.01	32.52	1686	0.02	91.24	1686	0.02	101.05							224.81
8/23/2010	1928	0.00	32.52	1928	0.00	91.24	1928	0.00	101.05							224.81
9/23/2010	2174	0.01	34.49	2174	0.02	96.27	2174	0.02	106.49							237.25
10/22/2010	2406	0.01	35.86	2406	0.01	98.85	2406	0.01	109.27							243.98
11/15/2010	2598	0.01	36.96	2598	0.01	101.41	2598	0.01	112.05							250.42
12/22/2010	2777	0.01	38.22	2955	0.02	107.99	2777	0.02	115.44							261.65
1/24/2011	2975	0.01	39.47	3352	0.01	110.39	2975	0.01	117.20							267.06
2/25/2011	3167	0.01	40.53	3737	0.01	114.08	3167	0.00	118.15							272.76
3/11/2011										222	1.72	381.87	218	0.35	75.54	730.17
3/18/2011	3293	0.01	41.27	3988	0.00	114.57	3293	0.00	118.34	366	0.51	453.50	362	0.20	104.77	832.46
3/25/2011										463	0.29	482.07	459	0.14	118.53	874.78
3/30/2011										558	0.32	512.25	553	0.08	126.48	912.92
4/8/2011										764	0.29	572.27	759	0.10	147.32	993.77
4/15/2011	3460	0.01	42.15	4322	0.00	115.07	3460	0.00	118.47	924	0.24	610.05	920	0.09	162.08	1047.81
5/19/2011	3665	0.00	42.87	4732	0.00	115.31	3665	0.00	118.53	1685	0.16	730.28	1681	0.09	233.92	1240.92
6/16/2011	3830	0.00	43.39	5062	0.00	115.55	3830	0.00	118.81	2191	0.11	753.86	2187	0.03	251.58	1283.20
7/15/2011	4472	0.00	44.96	4472	0.00	115.18	4472	0.00	119.39	2750	0.08	830.85	2745	0.03	269.61	1380.36
8/22/2011	4775	0.00	45.59	4775	0.00	115.40	4775	0.01	121.30	3133	0.10	868.97	3129	0.03	280.03	1431.44
9/15/2011	4968	0.00	45.93	4968	0.00	115.51	4968	0.00	121.91	3630	0.08	906.88	3626	0.01	287.36	1477.64
10/14/2011	5199	0.00	46.20	5199	0.00	115.57	5199	0.00	122.54	4226	0.05	935.35	4222	0.01	293.51	1513.18
11/21/2011	5503	0.00	46.43	5503	0.00	115.62	5503	0.00	123.00	5019	0.04	966.50	5015	0.01	298.43	1549.98
12/14/2011	5670	0.00	46.53	5670	0.00	115.65	5670	0.00	123.67	5343	0.03	975.34	5339	0.01	300.62	1561.80
1/19/2012	5974	0.00	46.69	5974	0.00	115.71	5974	0.00	124.59	5993	0.00	975.34	5958	0.00	300.62	1562.94
2/15/2012	6189	0.00	46.80	6189	0.00	115.74	6189	0.01	126.03	6368	0.03	986.48	6364	0.00	300.62	1575.67
3/15/2012	6421	0.00	46.89	6421	0.00	115.79	6421	0.01	127.43	6946	0.03	1005.89	6942	0.00	300.62	1596.62
4/19/2012	6701	0.00	47.04	6701	0.00	115.84	6701	0.00	128.02	7629	0.05	1038.74	7625	0.00	301.65	1631.30
5/16/2012	6916	0.00	47.18	6916	0.00	115.88	6916	0.00	128.27	8143	0.04	1060.30	8138	0.00	303.65	1655.28

Table 4.6
Mass Removal - Phase 1 and Phase 2 AS/SVE Systems
December 2009 - September 2015
UTC Aerospace Systems
Plants 1/2 Facility
Rockford, Illinois

Date	Cell 1			Cell 2			Cell 3			Cell 4			Cell 5			Total Cumulative Mass Removal (lb)
	Total Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	Run Time (hr)	Mass Removal Rate (lb/hr)	Cumulative Mass Removal (lb)	
Pulse -off period June 1, 2012 to August 14, 2012																
8/14/2012	7094	0.00	47.54	7094	0.00	116.20	7094	0.00	129.03	8546	0.05	1081.05	8541	0.01	306.52	1680.34
9/17/2012	7317	0.00	47.99	7317	0.00	116.40	7317	0.02	133.04	9033	0.04	1102.58	9029	0.01	311.31	1711.33
Pulse -off period September 17, 2012 to November 14, 2012																
11/15/2012	7320	0.00	48.00	7320	0.00	116.40	7320	0.00	133.05	9037	0.05	1102.78	9033	0.01	311.34	1711.56
12/14/2012	7518	0.00	48.24	7518	0.00	116.86	7518	0.00	133.94	9439	0.00	1103.57	9436	0.01	313.67	1716.27
Pulse -off period December 14, 2012 to February 26, 2013																
2/26/2013	7518	0.00	48.19	7518	0.00	116.86	7519	0.00	133.94	9439	0.00	1103.57	9511	0.00	313.72	1716.32
4/11/2013	7723	0.00	48.32	7723	0.00	116.97	8134	0.00	134.40	9876	0.00	1105.48	9952	0.02	322.58	1727.74
Pulse -off period April 11, 2013 to May 10, 2013																
5/10/2013	7724	0.00	48.32	7724	0.00	116.97	8135	0.00	134.40	9882	0.00	1105.50	9958	0.01	322.66	1727.85
7/15/2013	8039	0.00	48.86	8039	0.00	117.21	9082	0.00	134.70	10907	0.00	1108.40	10984	0.02	339.59	1748.76
Pulse -off period July 15, 2013 to September 9, 2013																
9/9/2013	8040	0.00	48.86	8040	0.00	117.21	9083	0.00	134.70	10914	0.00	1108.44	10991	0.01	339.65	1748.86
11/18/2013	8372	0.00	49.15	8372	0.00	117.30	10081	0.00	136.08	11992	0.00	1110.90	12069	0.02	356.69	1770.12
Pulse -off period November 18, 2013 to January 15, 2014																
1/15/2014	8651	0.00	49.36	8651	0.00	117.51	10916	0.00	136.88	11997	0.00	1110.91	12074	0.01	356.73	1771.39
3/14/2014	8894	0.00	49.48	8894	0.00	117.52	11645	0.00	137.13	12980	0.00	1112.65	13057	0.01	368.96	1785.75
Pulse -off period March 14, 2014 to May 15, 2014																
5/15/2014	8990	0.00	49.54	8990	0.00	117.64	11934	0.00	137.98	12986	0.00	1112.67	13063	0.01	369.01	1786.83
7/23/2014	9321	0.00	50.01	9321	0.00	117.79	12926	0.00	138.52	14627	0.00	1113.02	14714	0.00	371.61	1790.95
Pulse -off period July 23, 2014 to September 16, 2014																
9/16/2014	9494	0.00	50.32	9494	0.00	118.05	13099	0.00	138.77	14628	0.00	1113.03	14715	0.00	371.61	1791.78
11/14/2014	9777	0.00	50.45	9777	0.00	118.12	13948	0.00	139.44	16008	0.00	1116.04	16095	0.00	377.77	1801.82
Pulse -off period November 14, 2014 to January 9, 2015																
1/9/2015	9778	0.00	50.45	9778	0.00	118.12	13949	0.00	139.44	16015	0.00	1116.05	16102	0.00	377.78	1801.84
3/13/2015	10045	0.00	50.56	10045	0.00	118.15	14750	0.00	140.07	17178	0.00	1117.32	17322	0.00	382.64	1808.74
Pulse -off period March 13, 2015 to May 15, 2015																
5/15/2015	10046	0.00	50.56	10046	0.00	118.15	14751	0.00	140.07	17186	0.00	1117.34	17329	0.00	382.66	1808.77
7/16/2015	10343	0.00	50.92	10343	0.00	118.25	15641	0.00	140.72	18436	0.00	1121.16	18578	0.00	384.37	1815.42
Pulse -off period July 16, 2015 to September 22, 2015																
9/22/2015	10343	0.00	50.92	10343	0.00	118.26	15641	0.00	140.72	18439	0.00	1121.16	18580	0.00	384.37	1815.43

FIGURES



FIGURES

Figures



AECOM

AECOM
4320 WINFIELD ROAD
WARRENVILLE, ILLINOIS 60555
PHONE: (630) 829-2464
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Facility Location Map
Area 9/10 Remedial Action
Southeast Rockford Groundwater
Contamination Superfund Site
Rockford, IL

FIGURE NUMBER

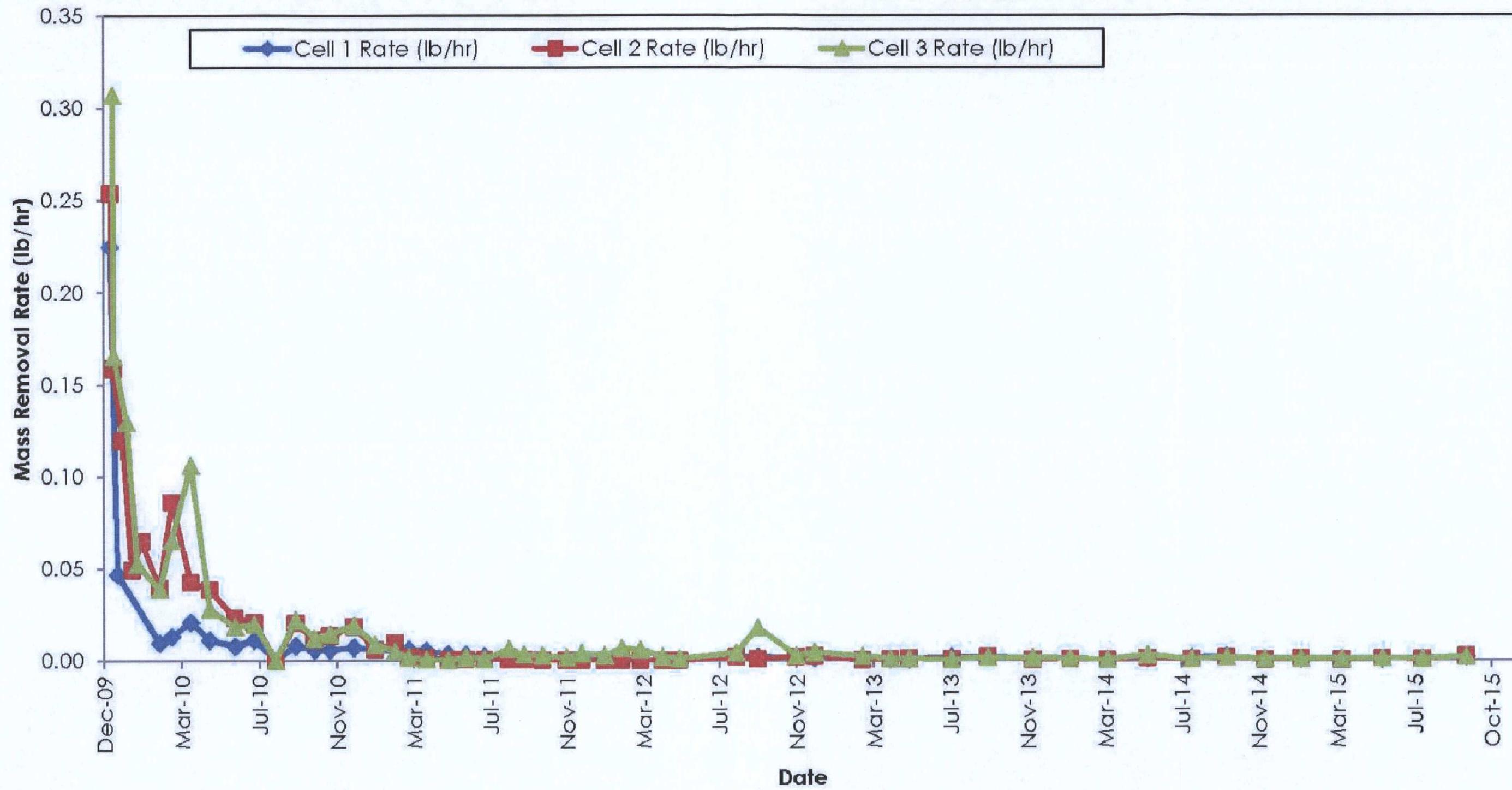
1

DRAWN BY:	DATE:	PROJECT NUMBER:	SHEET NUMBER:
JG	10/30/15	60339110.4214	1 of 1









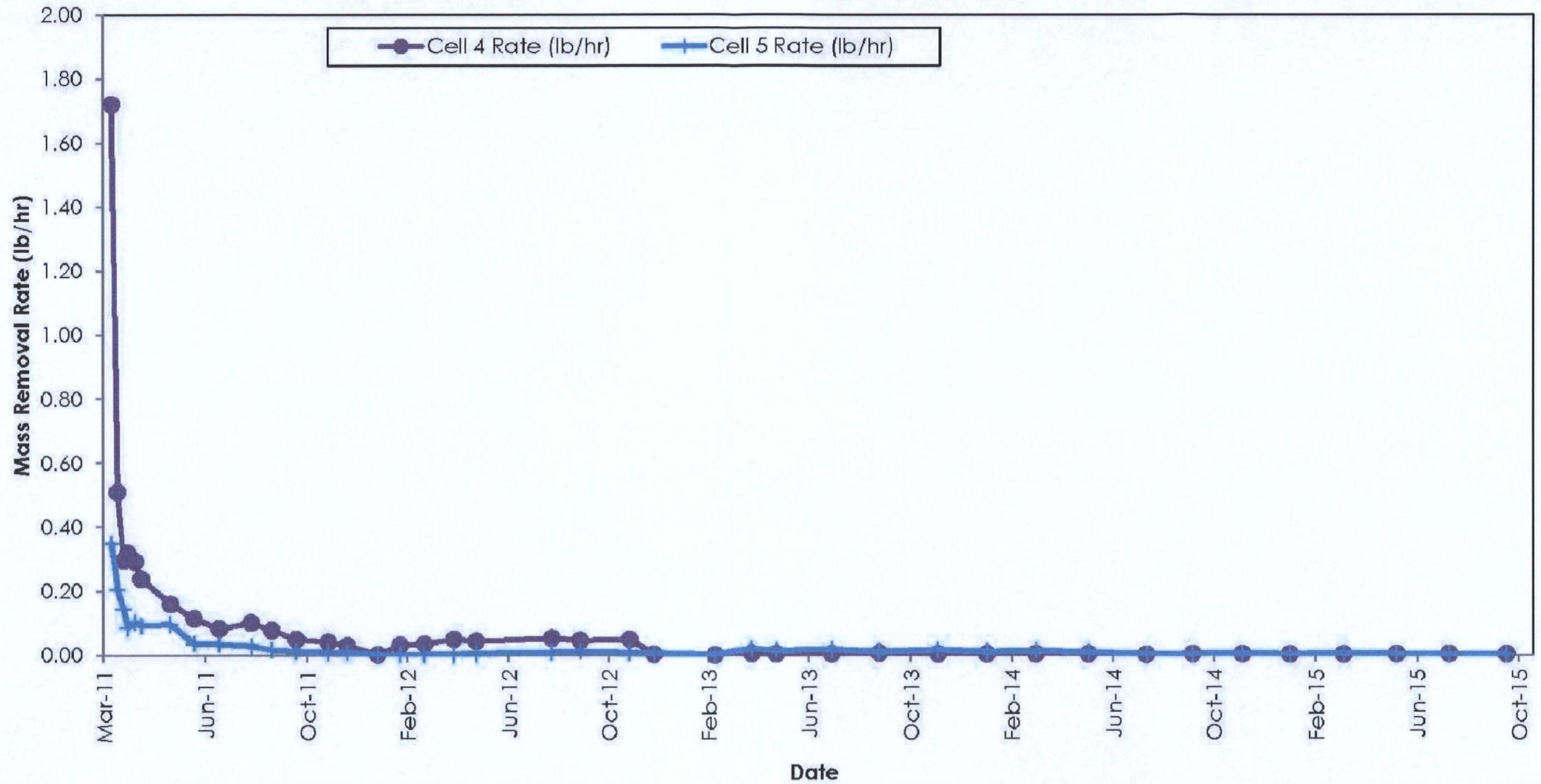
AECOM

AREA 9/10 REMEDIAL ACTION
ROCKFORD, ILLINOIS
PROJECT NO. 60339110.4214

AVERAGE VOC MASS REMOVAL RATE
VS TIME PHASE 1 AS/SVE SYSTEM

DATE: 10/30/15 DRWN: JG

FIGURE 7



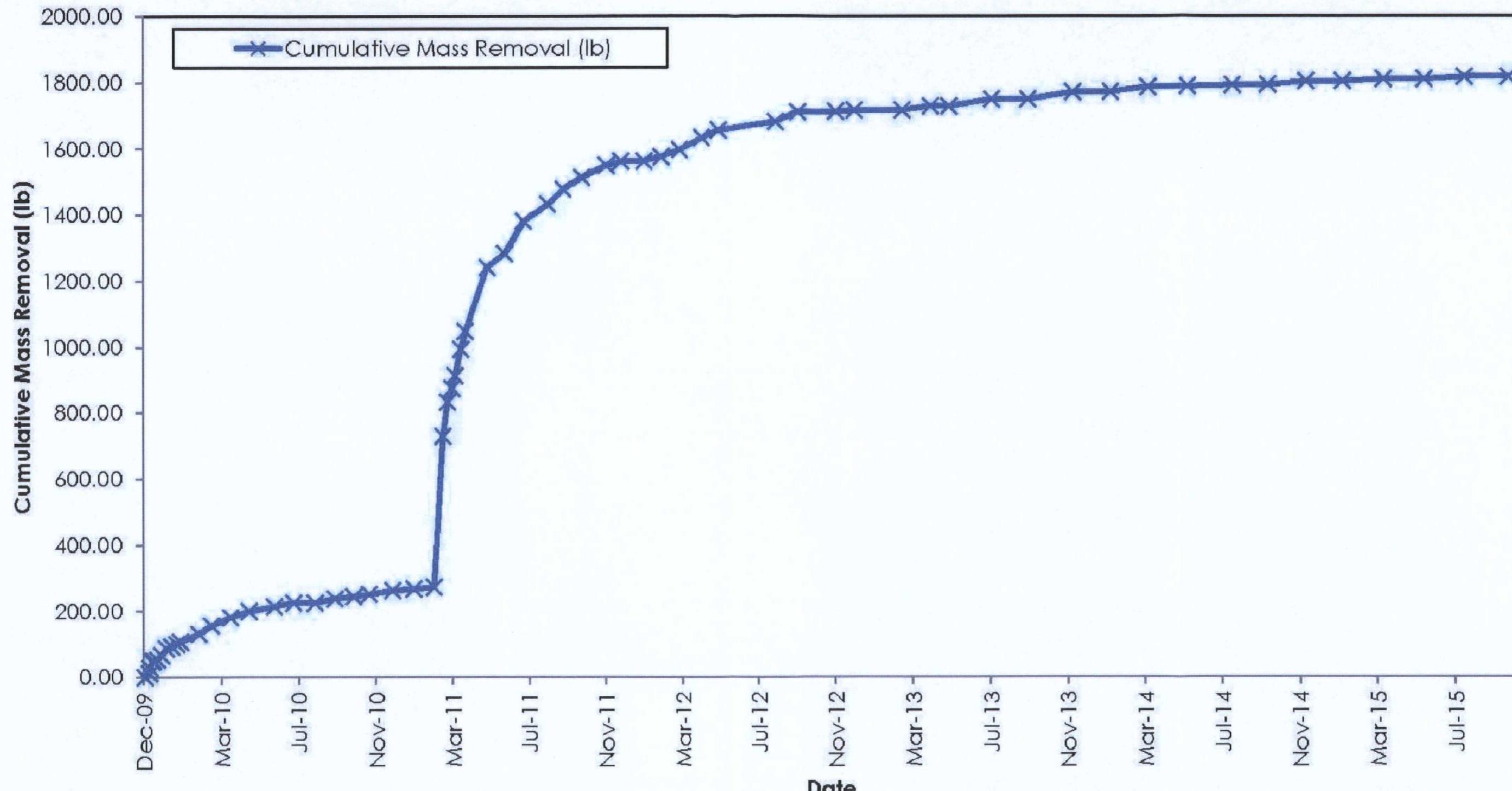
AECOM

AREA 9/10 REMEDIAL ACTION
ROCKFORD, ILLINOIS
PROJECT NO. 60339110.4214

AVERAGE VOC MASS REMOVAL RATE
VS TIME PHASE 2 AS/SVE SYSTEM

DATE: 10/30/15 DRWN: JG

FIGURE 8



AECOM

AREA 9/10 REMEDIAL ACTION
ROCKFORD, ILLINOIS
PROJECT NO. 60339110.4214

CUMULATIVE MASS REMOVAL
PHASE 1/ PHASE 2 AS/SVE SYSTEM

DATE: 10/30/15 DRWN: JG

FIGURE 9



A

!



A



Appendix A

Third Quarter 2015 GMZ and Performance Monitoring Well Analytical Data



Reissue #1
08/26/15

Technical Report for

United Technologies Corporation

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

60339110 PO#59048ACM

Accutest Job Number: JC1106

Sampling Dates: 08/04/15 - 08/06/15

Report to:

AECOM, INC.
27755 Diehl Road Suite 100
Warrenville, IL 60555
peter.hollatz@aecom.com

ATTN: Peter Hollatz

Total number of pages in report: 276



Test results contained within this data package meet the requirements
of the National Environmental Laboratory Accreditation Program
and/or state specific certification programs as applicable.

Nancy T. Cole

Nancy Cole
Laboratory Director

Client Service contact: Marie Meidhof 732-329-0200

Certifications: NJ(12129), NY(10983), CA, CT, DE, FL, IL, IN, KS, KY, LA, MA, MD, MI, MT, NC,
OH VAP (CL0056), AK (UST-103), AZ (AZ0786), PA, RI, SC, TN, TX, VA, WV, DoD ELAP (L-A-B L2248)

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Test results relate only to samples analyzed.



August 26, 2015

Mr. Peter Hollatz
AECOM
27755 Diehl Road, Suite 100
Warrenville, IL 60555

RE: Accutest job JC1106 - Reissue

Dear Mr. Hollatz,

The final report for Accutest job number JC1106 has been edited to reflect changes to your data package. These edits have been incorporated into the revised report which is attached.

The field ID for JC1106-9 has been revised to match the sample chain-of-custody. This report incorporates these revision.

Our apologies for any inconvenience the above issue may have caused you. Please contact me at 732-355-4552 if I may be of further assistance in this matter, or if you have any further questions regarding this data report.

Sincerely,

Accutest Laboratories

A handwritten signature in black ink that appears to read "Marie Meidhof".

Marie Meidhof
Project Manager

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Sample Summary

United Technologies Corporation

Job No: JC1106

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Collected Date	Time By	Received	Matrix Code	Type	Client Sample ID
JC1106-1	08/04/15	12:20 AH	08/08/15	AQ	Ground Water	HSSER-SMW08-080415
JC1106-2	08/04/15	13:35 AH	08/08/15	AQ	Ground Water	HSSER-GMZ01-080415
JC1106-3	08/04/15	14:25 AH	08/08/15	AQ	Ground Water	HSSER-SMW01-080415
JC1106-4	08/04/15	15:30 AH	08/08/15	AQ	Ground Water	HSSER-SMW02-080415
JC1106-5	08/04/15	16:55 AH	08/08/15	AQ	Ground Water	HSSER-MW203-080415
JC1106-6	08/05/15	08:25 AH	08/08/15	AQ	Ground Water	HSSER-MW07FGA-080515
JC1106-7	08/04/15	11:30 AH	08/08/15	AQ	Field Blank Water	HSSER-FBLK01-080415
JC1106-8	08/05/15	09:25 AH	08/08/15	AQ	Ground Water	HSSER-SMW19-080515
JC1106-9	08/05/15	10:55 AH	08/08/15	AQ	Ground Water	HSSER-GMZ04-080515
JC1106-10	08/05/15	12:00 AH	08/08/15	AQ	Ground Water	HSSER-SMW21-080515
JC1106-11	08/05/15	13:45 AH	08/08/15	AQ	Ground Water	HSSER-SMW20-080515
JC1106-12	08/05/15	15:00 AH	08/08/15	AQ	Ground Water	HSSER-GMZ03-080515
JC1106-13	08/05/15	16:30 AH	08/08/15	AQ	Ground Water	HSSER-GMZ02-080515

Sample Summary

(continued)

United Technologies Corporation

Job No: JC1106

ENSRLW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Collected Date	Time By	Received	Matrix Code Type	Client Sample ID
JC1106-13D	08/05/15	16:30 AH	08/08/15	AQ Water Dup/MSD	HSSER-MSD01-080515
JC1106-13S	08/05/15	16:30 AH	08/08/15	AQ Water Matrix Spike	HSSER-MS01-080515
JC1106-14	08/05/15	00:00 AH	08/08/15	AQ Ground Water	HSSER-DUP01-080515
JC1106-15	08/06/15	08:20 AH	08/08/15	AQ Ground Water	HSSER-PMW01-080615
JC1106-16	08/06/15	09:25 AH	08/08/15	AQ Ground Water	HSSER-PMW02-080615
JC1106-17	08/06/15	08:40 AH	08/08/15	AQ Equipment Blank	HSSER-EBLK01-080615
JC1106-18	08/06/15	10:30 AH	08/08/15	AQ Ground Water	HSSER-SMW04-080615
JC1106-19	08/06/15	10:30 AH	08/08/15	AQ Trip Blank Water	HSSER-TRIP01-080415



CASE NARRATIVE / CONFORMANCE SUMMARY

Client: United Technologies Corporation

Job No JC1106

Site: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Report Date 8/24/2015 11:18:11 A

On 08/08/2015, 16 Sample(s), 1 Trip Blank(s) and 1 Field Blank(s) and 1 Equipment Blank(s) were received at Accutest Laboratories at a maximum corrected temperature of 1.9 C. Samples were intact and chemically preserved, unless noted below. An Accutest Job Number of JC1106 was assigned to the project. Laboratory sample ID, client sample ID and dates of sample collection are detailed in the report's Results Summary Section.

Specified quality control criteria were achieved for this job except as noted below. For more information, please refer to the analytical results and QC summary pages.

Volatiles by GCMS By Method SW846 8260C

Matrix: AQ

Batch ID: V3D4822

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JC1106-13MS, JC1106-13MSD were used as the QC samples indicated.

Matrix: AQ

Batch ID: V3D4823

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JC1072-1MS, JC1072-1MSD were used as the QC samples indicated.

Matrix: AQ

Batch ID: V3D4824

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JC1106-9MS, JC1106-9MSD were used as the QC samples indicated.
- Matrix Spike Recovery(s) for 1,1,1-Trichloroethane are outside control limits. Outside control limits due to matrix interference.
- Matrix Spike Duplicate Recovery(s) for 1,1,1-Trichloroethane are outside control limits. Outside control limits due to matrix interference.

Accutest certifies that data reported for samples received, listed on the associated custody chain or analytical task order, were produced to specifications meeting Accutest's Quality System precision, accuracy and completeness objectives except as noted.

Estimated non-standard method measurement uncertainty data is available on request, based on quality control bias and implicit for standard methods. Acceptable uncertainty requires tested parameter quality control data to meet method criteria.

Accutest Laboratories is not responsible for data quality assumptions if partial reports are used and recommends that this report be used in its entirety. Data release is authorized by Accutest Laboratories indicated via signature on the report cover

Summary of Hits

Job Number: JC1106

Account: United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Collected: 08/04/15 thru 08/06/15



Lab Sample ID Analyte	Client Sample ID	Result/ Qual	RL	MDL	Units	Method
JC1106-1 HSSER-SMW08-080415						
1,1-Dichloroethane	0.0165	0.0010	0.00017	mg/l	SW846 8260C	
1,1-Dichloroethene	0.0011	0.0010	0.00051	mg/l	SW846 8260C	
cis-1,2-Dichloroethene	0.140	0.0010	0.00027	mg/l	SW846 8260C	
trans-1,2-Dichloroethene	0.0013	0.0010	0.00065	mg/l	SW846 8260C	
Tetrachloroethene	0.0243	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0040	0.0010	0.00025	mg/l	SW846 8260C	
Trichloroethene	0.0021	0.0010	0.00022	mg/l	SW846 8260C	
JC1106-2 HSSER-GMZ01-080415						
1,1-Dichloroethane	0.0066	0.0010	0.00017	mg/l	SW846 8260C	
cis-1,2-Dichloroethene	0.0037	0.0010	0.00027	mg/l	SW846 8260C	
Tetrachloroethene	0.0183	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0064	0.0010	0.00025	mg/l	SW846 8260C	
Trichloroethene	0.0018	0.0010	0.00022	mg/l	SW846 8260C	
JC1106-3 HSSER-SMW01-080415						
Tetrachloroethene	0.0013	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0013	0.0010	0.00025	mg/l	SW846 8260C	
JC1106-4 HSSER-SMW02-080415						
1,1-Dichloroethane	0.0011	0.0010	0.00017	mg/l	SW846 8260C	
cis-1,2-Dichloroethene	0.0014	0.0010	0.00027	mg/l	SW846 8260C	
Tetrachloroethene	0.00061 J	0.0010	0.00040	mg/l	SW846 8260C	
JC1106-5 HSSER-MW203-080415						
Tetrachloroethene	0.0037	0.0010	0.00040	mg/l	SW846 8260C	
JC1106-6 HSSER-MW07FGA-080515						
Tetrachloroethene	0.0011	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0011	0.0010	0.00025	mg/l	SW846 8260C	
JC1106-7 HSSER-FBLK01-080415						
Toluene	0.00045 J	0.0010	0.00016	mg/l	SW846 8260C	
JC1106-8 HSSER-SMW19-080515						
cis-1,2-Dichloroethene	0.00070 J	0.0010	0.00027	mg/l	SW846 8260C	

Summary of Hits

Job Number: JC1106

Account: United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Collected: 08/04/15 thru 08/06/15



Lab Sample ID Analyte	Client Sample ID	Result/ Qual	RL	MDL	Units	Method
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Tetrachloroethene	0.00086 J	0.0010	0.00040	mg/l	SW846 8260C
1,1,1-Trichloroethane	0.00034 J	0.0010	0.00025	mg/l	SW846 8260C
Trichloroethene	0.0106	0.0010	0.00022	mg/l	SW846 8260C

JC1106-9 HSSER-GMZ04-080515

1,1-Dichloroethane	0.0011	0.0010	0.00017	mg/l	SW846 8260C
1,1-Dichloroethene	0.00093 J	0.0010	0.00051	mg/l	SW846 8260C
cis-1,2-Dichloroethene	0.0035	0.0010	0.00027	mg/l	SW846 8260C
1,1,1-Trichloroethane	0.0536	0.0010	0.00025	mg/l	SW846 8260C
Trichloroethene	0.0026	0.0010	0.00022	mg/l	SW846 8260C

JC1106-10 HSSER-SMW21-080515

1,1-Dichloroethane	0.00025 J	0.0010	0.00017	mg/l	SW846 8260C
cis-1,2-Dichloroethene	0.00099 J	0.0010	0.00027	mg/l	SW846 8260C
1,1,1-Trichloroethane	0.010	0.0010	0.00025	mg/l	SW846 8260C
Trichloroethene	0.00084 J	0.0010	0.00022	mg/l	SW846 8260C

JC1106-11 HSSER-SMW20-080515

No hits reported in this sample.

JC1106-12 HSSER-GMZ03-080515

1,1-Dichloroethane	0.00028 J	0.0010	0.00017	mg/l	SW846 8260C
1,1,1-Trichloroethane	0.00035 J	0.0010	0.00025	mg/l	SW846 8260C

JC1106-13 HSSER-GMZ02-080515

1,1-Dichloroethane	0.0016	0.0010	0.00017	mg/l	SW846 8260C
cis-1,2-Dichloroethene	0.00057 J	0.0010	0.00027	mg/l	SW846 8260C
Tetrachloroethene	0.00043 J	0.0010	0.00040	mg/l	SW846 8260C
1,1,1-Trichloroethane	0.0030	0.0010	0.00025	mg/l	SW846 8260C
Trichloroethene	0.00040 J	0.0010	0.00022	mg/l	SW846 8260C

JC1106-14 HSSER-DUP01-080515

1,1-Dichloroethane	0.00026 J	0.0010	0.00017	mg/l	SW846 8260C
1,1,1-Trichloroethane	0.00036 J	0.0010	0.00025	mg/l	SW846 8260C

JC1106-15 HSSER-PMW01-080615

1,1-Dichloroethane	0.0086	0.0010	0.00017	mg/l	SW846 8260C
1,1-Dichloroethene	0.0018	0.0010	0.00051	mg/l	SW846 8260C

Summary of Hits

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Job Number: JC1106

Account: United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Collected: 08/04/15 thru 08/06/15



Lab Sample ID	Client Sample ID	Result/ Qual	RL	MDL	Units	Method
cis-1,2-Dichloroethene	0.0011	0.0010	0.00027	mg/l	SW846 8260C	
Tetrachloroethene	0.0192	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0547	0.0010	0.00025	mg/l	SW846 8260C	
Trichloroethene	0.0013	0.0010	0.00022	mg/l	SW846 8260C	
JC1106-16 HSSER-PMW02-080615						
1,1-Dichloroethane	0.0069	0.0010	0.00017	mg/l	SW846 8260C	
cis-1,2-Dichloroethene	0.0043	0.0010	0.00027	mg/l	SW846 8260C	
Tetrachloroethene	0.0088	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0042	0.0010	0.00025	mg/l	SW846 8260C	
Trichloroethene	0.0018	0.0010	0.00022	mg/l	SW846 8260C	
Vinyl chloride	0.0016	0.0010	0.00015	mg/l	SW846 8260C	
JC1106-17 HSSER-EBLK01-080615						
Toluene	0.00033 J	0.0010	0.00016	mg/l	SW846 8260C	
JC1106-18 HSSER-SMW04-080615						
1,1-Dichloroethane	0.0121	0.0010	0.00017	mg/l	SW846 8260C	
cis-1,2-Dichloroethene	0.0756	0.0010	0.00027	mg/l	SW846 8260C	
Tetrachloroethene	0.0090	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0031	0.0010	0.00025	mg/l	SW846 8260C	
Trichloroethene	0.0023	0.0010	0.00022	mg/l	SW846 8260C	
Vinyl chloride	0.0163	0.0010	0.00015	mg/l	SW846 8260C	
JC1106-19 HSSER-TRIP01-080415						
No hits reported in this sample.						

No hits reported in this sample.



Sample Results

Report of Analysis

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-SMW08-080415
Lab Sample ID: JC1106-1
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/04/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112575.D	1	08/12/15	XC	n/a	n/a	V3D4824
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0165	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	0.0011	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.140	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	0.0013	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0243	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0040	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.0021	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	112%		76-120%
17060-07-0	1,2-Dichloroethane-D4	110%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-GMZ01-080415
Lab Sample ID: JC1106-2
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/04/15
Date Received: 08/08/15
Percent Solids: n/a

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112574.D	1	08/12/15	XC	n/a	n/a	V3D4824
Run #2							

Purge Volume	
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0066	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.0037	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0183	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0064	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.0018	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	112%		76-120%
17060-07-0	1,2-Dichloroethane-D4	115%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	95%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-SMW01-080415**Lab Sample ID:** JC1106-3**Date Sampled:** 08/04/15**Matrix:** AQ - Ground Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112531.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

Purge Volume

Run #1 5.0 ml

Run #2

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0013	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0013	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	110%		76-120%
17060-07-0	1,2-Dichloroethane-D4	113%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	98%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-SMW02-080415
Lab Sample ID: JC1106-4
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/04/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112532.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0011	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.0014	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.00061	0.0010	0.00040	mg/l	J
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	114%		76-120%
17060-07-0	1,2-Dichloroethane-D4	113%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-MW203-080415
Lab Sample ID: JC1106-5
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/04/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112523.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

Purge Volume
Run #1 5.0 ml
Run #2

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0037	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	112%		76-120%
17060-07-0	1,2-Dichloroethane-D4	114%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected MDL = Method Detection Limit
RL = Reporting Limit
E = Indicates value exceeds calibration range

J = Indicates an estimated value
B = Indicates analyte found in associated method blank
N = Indicates presumptive evidence of a compound

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Client Sample ID: HSSER-MW07FGA-080515**Lab Sample ID:** JC1106-6**Date Sampled:** 08/05/15**Matrix:** AQ - Ground Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112530.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

Purge Volume

Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0011	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0011	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	111%		76-120%
17060-07-0	1,2-Dichloroethane-D4	112%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	96%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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Client Sample ID: HSSER-FBLK01-080415
Lab Sample ID: JC1106-7
Matrix: AQ - Field Blank Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/04/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112522.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	0.00045	0.0010	0.00016	mg/l	J
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	111%		76-120%
17060-07-0	1,2-Dichloroethane-D4	112%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	96%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

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**Client Sample ID:** HSSER-SMW19-080515**Lab Sample ID:** JC1106-8**Matrix:** AQ - Ground Water**Method:** SW846 8260C**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL**Date Sampled:** 08/05/15**Date Received:** 08/08/15**Percent Solids:** n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112535.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.00070	0.0010	0.00027	mg/l	J
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.00086	0.0010	0.00040	mg/l	J
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.00034	0.0010	0.00025	mg/l	J
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.0106	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	113%		76-120%
17060-07-0	1,2-Dichloroethane-D4	115%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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Client Sample ID: HSSER-GMZ04-080515**Lab Sample ID:** JC1106-9**Matrix:** AQ - Ground Water**Method:** SW846 8260C**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL**Date Sampled:** 08/05/15**Date Received:** 08/08/15**Percent Solids:** n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112572.D	1	08/12/15	XC	n/a	n/a	V3D4824
Run #2							

Purge Volume

Run #1 5.0 ml

Run #2

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0011	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	0.00093	0.0010	0.00051	mg/l	J
156-59-2	cis-1,2-Dichloroethene	0.0035	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0536	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.0026	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	112%		76-120%
17060-07-0	1,2-Dichloroethane-D4	111%		73-122%
2037-26-5	Toluene-D8	102%		84-119%
460-00-4	4-Bromofluorobenzene	96%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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Client Sample ID: HSSER-SMW21-080515
Lab Sample ID: JC1106-10
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/05/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112525.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.00025	0.0010	0.00017	mg/l	J
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.00099	0.0010	0.00027	mg/l	J
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.010	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.00084	0.0010	0.00022	mg/l	J
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	113%		76-120%
17060-07-0	1,2-Dichloroethane-D4	114%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

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Client Sample ID: HSSER-SMW20-080515
Lab Sample ID: JC1106-11
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/05/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112538.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	114%		76-120%
17060-07-0	1,2-Dichloroethane-D4	113%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

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4**Client Sample ID:** HSSER-GMZ03-080515**Lab Sample ID:** JC1106-12**Matrix:** AQ - Ground Water**Method:** SW846 8260C**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL**Date Sampled:** 08/05/15**Date Received:** 08/08/15**Percent Solids:** n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112537.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

Purge Volume

Run #1 5.0 ml

Run #2

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.00028	0.0010	0.00017	mg/l	J
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.00035	0.0010	0.00025	mg/l	J
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	113%		76-120%
17060-07-0	1,2-Dichloroethane-D4	115%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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Client Sample ID: HSSER-GMZ02-080515

Lab Sample ID: JC1106-13

Date Sampled: 08/05/15

Matrix: AQ - Ground Water

Date Received: 08/08/15

Method: SW846 8260C

Percent Solids: n/a

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112524.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

Purge Volume

Run #1 5.0 ml

Run #2

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0016	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.00057	0.0010	0.00027	mg/l	J
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.00043	0.0010	0.00040	mg/l	J
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0030	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.00040	0.0010	0.00022	mg/l	J
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	111%		76-120%
17060-07-0	1,2-Dichloroethane-D4	112%		73-122%
2037-26-5	Toluene-D8	98%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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Client Sample ID:	HSSER-DUP01-080515	Date Sampled:	08/05/15
Lab Sample ID:	JC1106-14	Date Received:	08/08/15
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Method:	SW846 8260C		
Project:	ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112536.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.00026	0.0010	0.00017	mg/l	J
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.00036	0.0010	0.00025	mg/l	J
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	114%		76-120%
17060-07-0	1,2-Dichloroethane-D4	117%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	97%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

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Client Sample ID: HSSER-PMW01-080615**Lab Sample ID:** JC1106-15**Date Sampled:** 08/06/15**Matrix:** AQ - Ground Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112533.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0086	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	0.0018	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.0011	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0192	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0547	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.0013	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	113%		76-120%
17060-07-0	1,2-Dichloroethane-D4	112%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	96%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID:	HSSER-PMW02-080615	Date Sampled:	08/06/15
Lab Sample ID:	JC1106-16	Date Received:	08/08/15
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Method:	SW846 8260C		
Project:	ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112573.D	1	08/12/15	XC	n/a	n/a	V3D4824
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0069	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.0043	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0088	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0042	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.0018	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	0.0016	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	111%		76-120%
17060-07-0	1,2-Dichloroethane-D4	112%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	95%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID:	HSSER-EBLK01-080615	Date Sampled:	08/06/15
Lab Sample ID:	JC1106-17	Date Received:	08/08/15
Matrix:	AQ - Equipment Blank	Percent Solids:	n/a
Method:	SW846 8260C		
Project:	ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112555.D	1	08/12/15	XC	n/a	n/a	V3D4823
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	0.00033	0.0010	0.00016	mg/l	J
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	108%		76-120%
17060-07-0	1,2-Dichloroethane-D4	112%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	98%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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4**Client Sample ID:** HSSER-SMW04-080615**Lab Sample ID:** JC1106-18**Date Sampled:** 08/06/15**Matrix:** AQ - Ground Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112534.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0121	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.0756	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0090	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0031	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.0023	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	0.0163	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	111%		76-120%
17060-07-0	1,2-Dichloroethane-D4	112%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	98%		78-117%

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: HSSER-TRIP01-080415**Lab Sample ID:** JC1106-19**Date Sampled:** 08/06/15**Matrix:** AQ - Trip Blank Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	3D112529.D	1	08/11/15	XC	n/a	n/a	V3D4822
Run #2							

Purge Volume

Run #1 5.0 ml

Run #2

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	112%		76-120%
17060-07-0	1,2-Dichloroethane-D4	113%		73-122%
2037-26-5	Toluene-D8	101%		84-119%
460-00-4	4-Bromofluorobenzene	96%		78-117%

ND = Not detected

MDL = Method Detection Limit

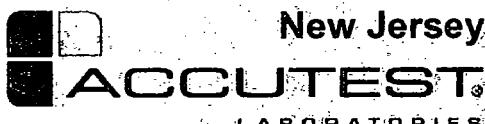
J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound



Misc. Forms



Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody
- Sample Tracking Chronicle
- Internal Chain of Custody



GW
WTB

CHAIN OF CUSTODY

2235 Route 130, Dayton, NJ 08810
TEL 732-329-0200 FAX: 732-329-3499/3480
www.acme1st.com

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JC1106: Chain of Custody

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bw
WTR

CHAIN OF CUSTODY

PAGE 2 OF 2

2235 Route 130, Dayton, NJ 08810
TEL: 732-329-0200 FAX: 732-329-3480
www.accutest.com

FED-EX Tracking #	Boiler Order Control #
6034940616410	Accutest Job #
Accutest Quote #	JC1106

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)												Matrix Codes	
Company Name AECOM	Project Name UTMS Plants 1/2 Fac J1.tz	Street	City	Billing Information (if different from Report to)												DW - Drinking Water	
Street Address 27755 Doh Rd Suite 100	State NY	City Wheatonville Fl	State FL	Company Name	Street Address	City	State	Zip	WW - Water								
City Wheatonville Fl	Zip 32055	City Moskford	State FL	Client Purchase Order # 10339110	City	State	Zip	SW - Surface Water									
Phone # (630-836-5366	Fax # 630-836-1711	Phone # 59048 ACM	Project Manager Peter Hallatz	Attention	SO - Sot	SL - Sludge	SED - Sediment	DI - DN									
Sampler(s) Name(s) Allen Hallatz	Phone # 630-836-1700	Project Manager Peter Hallatz			LIQ - Oil & Liquid	AR - Ash	SOL - Other Solid	WP - Wipe									
					ED - Equipment Blank	RB - Rinse Blank	TB - Trip Blank										
Collection																LAB USE ONLY	
Method Sample #		Field ID / Point of Collection	MECH/DIV ID #	Date	Time	Sampled by	Matrix	# of bottles	HC	NC	PC	PA	ME	MD	ENH		
									X	X	X	X	X	X	X		
		HSSER-6M202-080515		8-5-15	1630	FX	GW	3									
		HSSER-1M101-080515		8-5-15	1630			3	X								
		HSSER-1M501-080515		8-5-15	1630			3	X								
		HSSER-DUPO1-080515		8-5-15	0000			3	X								
		HSSER-PML01-080615		X	0820			3	X								
		HSSER-PML02-080615		8-6-15	0925			3	X								
		HSSER-EBLK01-080615		8-6-15	0840			3	X								
		HSSER-B-SMG04-080615		8-6-15	1030			3	X								
		HSSER-TMP01-080415		8-4-15	—			2	X								
Data Deliverable Information																Comments / Special Instructions	
Turnaround Time (Business days)																List of 13 VOC	
Approved By (Accutest PM) / Date:								NYASP Category A									
<input type="checkbox"/> Std. 10 Business Days								<input type="checkbox"/> Commercial "A" (Level 1)									
<input type="checkbox"/> 5 Day RUSH								<input type="checkbox"/> NYASP Category B									
<input type="checkbox"/> 3 Day RUSH								<input type="checkbox"/> Commercial "B" (Level 2)									
<input type="checkbox"/> 2 Day RUSH								<input type="checkbox"/> FULLY1 (Level 3+4)									
<input type="checkbox"/> 1 Day RUSH								<input type="checkbox"/> State Forms									
<input type="checkbox"/> other								<input type="checkbox"/> EDD Format									
Emergency & RUSH T/A date available VIA Lablink								<input type="checkbox"/> Other _____									
NJ Data of Known Quality Protocol Reporting								NJ Data _____									
Commercial "A" = Results Only, Commercial "B" = Results + QC Summary								Commercial "C" = Results + QC Summary + Partial Raw data									
NJ Reduced = Results + QC Summary + Partial Raw data								List of 13 VOC									
Sample Custody must be documented below each time samples change possession, including courier delivery.																Level 1C Data	
Retainquished by Sampler: 1 Allen Hallatz		Date Time: 8-7-15 1800	Received By: 1 FX	Retainquished By: 2		Date Time: 8-6-15 18:00	Received By: 2										
Retainquished by Sampler: 3		Date Time: —	Received By: 3	Retainquished By: 4		Date Time: —	Received By: 4										
Retainquished by: 5		Date Time: —	Received By: 5	Custody Seal #		<input type="checkbox"/> intact	Preserved where applicable	<input type="checkbox"/>	On Ice	Cooler Temp	24c	-ip					

JC1106: Chain of Custody
Page 2 of 3



Accutest Laboratories Sample Receipt Summary

Accutest Job Number: JC1106

Client: _____

Project: _____

Date / Time Received: 8/8/2015 10:00:00 AM

Delivery Method: _____

Airbill #'s: _____

Cooler Temps (Raw Measured) °C: Cooler 1: (2.2);

Cooler Temps (Corrected) °C: Cooler 1: (1.9);

Cooler Security Y or N

- | | | | | | |
|---------------------------|-------------------------------------|--------------------------|-----------------------|-------------------------------------|--------------------------|
| 1. Custody Seals Present: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 3. COC Present: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Custody Seals Intact: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | 4. Smpl Dates/Time OK | <input checked="" type="checkbox"/> | <input type="checkbox"/> |

Cooler Temperature Y or N

- | | | |
|------------------------------|-------------------------------------|--------------------------|
| 1. Temp criteria achieved: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Cooler temp verification: | IR Gun | |
| 3. Cooler media: | Ice (Bag) | |
| 4. No. Coolers: | 1 | |

Quality Control Preservation Y or N N/A

- | | | | |
|---------------------------------|-------------------------------------|--------------------------|--------------------------|
| 1. Trip Blank present / cooler: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 2. Trip Blank listed on COC: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |
| 3. Samples preserved properly: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | |
| 4. VOCs headspace free: | <input checked="" type="checkbox"/> | <input type="checkbox"/> | <input type="checkbox"/> |

Sample Integrity - Documentation

- | | | |
|--|-------------------------------------|--------------------------|
| 1. Sample labels present on bottles: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Container labeling complete: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 3. Sample container label / COC agree: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |

Sample Integrity - Condition

- | | | |
|----------------------------------|-------------------------------------|--------------------------|
| 1. Sample recvd within HT: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. All containers accounted for: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 3. Condition of sample: | Intact | |

Sample Integrity - Instructions

- | | | |
|---|-------------------------------------|-------------------------------------|
| 1. Analysis requested is clear: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 2. Bottles received for unspecified tests | <input type="checkbox"/> | <input checked="" type="checkbox"/> |
| 3. Sufficient volume recvd for analysis: | <input checked="" type="checkbox"/> | <input type="checkbox"/> |
| 4. Compositing instructions clear: | <input type="checkbox"/> | <input type="checkbox"/> |
| 5. Filtering instructions clear: | <input type="checkbox"/> | <input checked="" type="checkbox"/> |

Comments

Accutest Laboratories
V:732.329.02002235 US Highway 130
F: 732.329.3499Dayton, New Jersey
www.accutest.com**JC1106: Chain of Custody****Page 3 of 3**

Internal Sample Tracking Chronicle

United Technologies Corporation

Job No: JC1106

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Method	Analyzed	By	Prepped	By	Test Codes
JC1106-1 HSSER-SMW08-080415	Collected: 04-AUG-15 12:20 By: AH			Received: 08-AUG-15 By: DG		
JC1106-1 SW846 8260C		12-AUG-15 14:48 XC				V8260SL
JC1106-2 HSSER-GMZ01-080415	Collected: 04-AUG-15 13:35 By: AH			Received: 08-AUG-15 By: DG		
JC1106-2 SW846 8260C		12-AUG-15 14:21 XC				V8260SL
JC1106-3 HSSER-SMW01-080415	Collected: 04-AUG-15 14:25 By: AH			Received: 08-AUG-15 By: DG		
JC1106-3 SW846 8260C		11-AUG-15 18:18 XC				V8260SL
JC1106-4 HSSER-SMW02-080415	Collected: 04-AUG-15 15:30 By: AH			Received: 08-AUG-15 By: DG		
JC1106-4 SW846 8260C		11-AUG-15 18:45 XC				V8260SL
JC1106-5 HSSER-MW203-080415	Collected: 04-AUG-15 16:55 By: AH			Received: 08-AUG-15 By: DG		
JC1106-5 SW846 8260C		11-AUG-15 14:36 XC				V8260SL
JC1106-6 HSSER-MW07FGA-080515	Collected: 05-AUG-15 08:25 By: AH			Received: 08-AUG-15 By: DG		
JC1106-6 SW846 8260C		11-AUG-15 17:51 XC				V8260SL
JC1106-7 HSSER-FBLK01-080415	Collected: 04-AUG-15 11:30 By: AH			Received: 08-AUG-15 By: DG		
JC1106-7 SW846 8260C		11-AUG-15 14:09 XC				V8260SL
JC1106-8 HSSER-SMW19-080515	Collected: 05-AUG-15 09:25 By: AH			Received: 08-AUG-15 By: DG		
JC1106-8 SW846 8260C		11-AUG-15 20:07 XC				V8260SL

Internal Sample Tracking Chronicle

United Technologies Corporation

Job No: JC1106

ENSRLW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Method	Analyzed	By	Prepped	By	Test Codes
JC1106-9	Collected: 05-AUG-15 10:55 By: AH HSSER-GMZ04-080515			Received: 08-AUG-15 By: DG		
JC1106-9	SW846 8260C	12-AUG-15 13:27 XC			V8260SL	
JC1106-10	Collected: 05-AUG-15 12:00 By: AH HSSER-SMW21-080515			Received: 08-AUG-15 By: DG		
JC1106-10	SW846 8260C	11-AUG-15 15:30 XC			V8260SL	
JC1106-11	Collected: 05-AUG-15 13:45 By: AH HSSER-SMW20-080515			Received: 08-AUG-15 By: DG		
JC1106-11	SW846 8260C	11-AUG-15 21:28 XC			V8260SL	
JC1106-12	Collected: 05-AUG-15 15:00 By: AH HSSER-GMZ03-080515			Received: 08-AUG-15 By: DG		
JC1106-12	SW846 8260C	11-AUG-15 21:01 XC			V8260SL	
JC1106-13	Collected: 05-AUG-15 16:30 By: AH HSSER-GMZ02-080515			Received: 08-AUG-15 By: DG		
JC1106-13	SW846 8260C	11-AUG-15 15:03 XC			V8260SL	
JC1106-14	Collected: 05-AUG-15 00:00 By: AH HSSER-DUP01-080515			Received: 08-AUG-15 By: DG		
JC1106-14	SW846 8260C	11-AUG-15 20:34 XC			V8260SL	
JC1106-15	Collected: 06-AUG-15 08:20 By: AH HSSER-PMW01-080615			Received: 08-AUG-15 By: DG		
JC1106-15	SW846 8260C	11-AUG-15 19:13 XC			V8260SL	
JC1106-16	Collected: 06-AUG-15 09:25 By: AH HSSER-PMW02-080615			Received: 08-AUG-15 By: DG		
JC1106-16	SW846 8260C	12-AUG-15 13:54 XC			V8260SL	

Internal Sample Tracking Chronicle

United Technologies Corporation

Job No: JC1106

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Method	Analyzed By	Prepped By	Test Codes
JC1106-17	Collected: 06-AUG-15 08:40 By: AH HSSER-EBLK01-080615		Received: 08-AUG-15 By: DG	
JC1106-17	SW846 8260C	12-AUG-15 05:07 XC		V8260SL
JC1106-18	Collected: 06-AUG-15 10:30 By: AH HSSER-SMW04-080615		Received: 08-AUG-15 By: DG	
JC1106-18	SW846 8260C	11-AUG-15 19:40 XC		V8260SL
JC1106-19	Collected: 06-AUG-15 10:30 By: AH HSSER-TRIP01-080415		Received: 08-AUG-15 By: DG	
JC1106-19	SW846 8260C	11-AUG-15 17:24 XC		V8260SL

Accutest Internal Chain of Custody

Page 1 of 3

Job Number: JC1106
Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Received: 08/08/15

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JC1106-1.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-1.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-1.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-1.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-1.2	Secured Storage	Ximena Collado	08/12/15 11:56	Retrieve from Storage
JC1106-1.2	Ximena Collado	GCMS3D	08/12/15 11:56	Load on Instrument
JC1106-1.2	GCMS3D	Ximena Collado	08/13/15 09:49	Unload from Instrument
JC1106-1.2	Ximena Collado	Secured Storage	08/13/15 09:49	Return to Storage
JC1106-2.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-2.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-2.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-2.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-2.2	Secured Storage	Ximena Collado	08/12/15 11:56	Retrieve from Storage
JC1106-2.2	Ximena Collado	GCMS3D	08/12/15 11:56	Load on Instrument
JC1106-2.2	GCMS3D	Ximena Collado	08/13/15 09:49	Unload from Instrument
JC1106-2.2	Ximena Collado	Secured Storage	08/13/15 09:49	Return to Storage
JC1106-3.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-3.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-3.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-3.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-4.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-4.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-4.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-4.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-5.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-5.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-5.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-5.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-6.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-6.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-6.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-6.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-7.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-7.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-7.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-7.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage

Accutest Internal Chain of Custody

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Job Number: JC1106
Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Received: 08/08/15

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JC1106-8.2	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-8.2	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-8.2	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-8.2	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-9.1	Secured Storage	Ximena Collado	08/12/15 11:56	Retrieve from Storage
JC1106-9.1	Ximena Collado	GCMS3D	08/12/15 11:56	Load on Instrument
JC1106-9.1	GCMS3D	Ximena Collado	08/13/15 09:49	Unload from Instrument
JC1106-9.1	Ximena Collado	Secured Storage	08/13/15 09:49	Return to Storage
JC1106-9.2	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-9.2	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-9.2	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-9.2	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-9.2	Secured Storage	Ximena Collado	08/12/15 16:36	Retrieve from Storage
JC1106-9.2	Ximena Collado	GCMS3D	08/12/15 16:36	Load on Instrument
JC1106-9.2	GCMS3D	Ximena Collado	08/13/15 09:49	Unload from Instrument
JC1106-9.2	Ximena Collado	Secured Storage	08/13/15 09:49	Return to Storage
JC1106-10.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-10.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-10.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-10.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-11.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-11.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-11.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-11.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-12.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-12.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-12.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-12.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-13.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-13.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-13.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-13.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-13.6	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-13.6	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-13.6	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-13.6	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage

Accutest Internal Chain of Custody

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Job Number: JC1106
Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Received: 08/08/15

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JC1106-13.7	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-13.7	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-13.7	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-13.7	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-14.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-14.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-14.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-14.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-15.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-15.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-15.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-15.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-16.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-16.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-16.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-16.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-16.2	Secured Storage	Ximena Collado	08/12/15 11:56	Retrieve from Storage
JC1106-16.2	Ximena Collado	GCMS3D	08/12/15 11:56	Load on Instrument
JC1106-16.2	GCMS3D	Ximena Collado	08/13/15 09:49	Unload from Instrument
JC1106-16.2	Ximena Collado	Secured Storage	08/13/15 09:49	Return to Storage
JC1106-17.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-17.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-17.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-17.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-18.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-18.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-18.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-18.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage
JC1106-19.1	Secured Storage	Ximena Collado	08/11/15 13:05	Retrieve from Storage
JC1106-19.1	Ximena Collado	GCMS3D	08/11/15 13:05	Load on Instrument
JC1106-19.1	GCMS3D	Ximena Collado	08/12/15 09:30	Unload from Instrument
JC1106-19.1	Ximena Collado	Secured Storage	08/12/15 09:30	Return to Storage



GC/MS Volatiles



QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries
- Instrument Performance Checks (BFB)
- Internal Standard Area Summaries
- Surrogate Recovery Summaries
- Initial and Continuing Calibration Summaries

Method Blank Summary**Job Number:** JC1106**Account:** UTC United Technologies Corporation**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3D4822-MB	3D112519.D	1	08/11/15	XC	n/a	n/a	V3D4822

The QC reported here applies to the following samples:**Method:** SW846 8260C

JC1106-3, JC1106-4, JC1106-5, JC1106-6, JC1106-7, JC1106-8, JC1106-10, JC1106-11, JC1106-12, JC1106-13, JC1106-14, JC1106-15, JC1106-18, JC1106-19

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	

CAS No. Surrogate Recoveries**Limits**

1868-53-7	Dibromofluoromethane	108%	76-120%
17060-07-0	1,2-Dichloroethane-D4	112%	73-122%
2037-26-5	Toluene-D8	102%	84-119%
460-00-4	4-Bromofluorobenzene	100%	78-117%

Method Blank Summary

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3D4823-MB	3D112544.D	1	08/12/15	XC	n/a	n/a	V3D4823

The QC reported here applies to the following samples:

Method: SW846 8260C

JC1106-17

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CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	

CAS No.	Surrogate Recoveries	Limits
1868-53-7	Dibromofluoromethane	112% 76-120%
17060-07-0	1,2-Dichloroethane-D4	114% 73-122%
2037-26-5	Toluene-D8	102% 84-119%
460-00-4	4-Bromofluorobenzene	96% 78-117%

CAS No.	Tentatively Identified Compounds	R.T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

Method Blank Summary**Job Number:** JC1106**Account:** UTC United Technologies Corporation**Project:** ENSRJLW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3D4824-MB	3D112569.D	1	08/12/15	XC	n/a	n/a	V3D4824

The QC reported here applies to the following samples:**Method:** SW846 8260C

JC1106-1, JC1106-2, JC1106-9, JC1106-16

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CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	

CAS No. Surrogate Recoveries Limits

1868-53-7	Dibromofluoromethane	114%	76-120%
17060-07-0	1,2-Dichloroethane-D4	110%	73-122%
2037-26-5	Toluene-D8	101%	84-119%
460-00-4	4-Bromofluorobenzene	97%	78-117%

CAS No. Tentatively Identified Compounds R.T. Est. Conc. Units Q

Total TIC, Volatile 0 ug/l

Method Blank Summary

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3D4823-MB2	3D112565.D	1	08/12/15	XC	n/a	n/a	V3D4823

The QC reported here applies to the following samples:

Method: SW846 8260C

JC1072-1MS, JC1072-1MSD

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	

CAS No. Surrogate Recoveries

	Limits
1868-53-7 Dibromofluoromethane	76-120%
17060-07-0 1,2-Dichloroethane-D4	73-122%
2037-26-5 Toluene-D8	84-119%
460-00-4 4-Bromofluorobenzene	78-117%

Blank Spike Summary

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3D4822-BS	3D112520.D	1	08/11/15	XC	n/a	n/a	V3D4822

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC1106-3, JC1106-4, JC1106-5, JC1106-6, JC1106-7, JC1106-8, JC1106-10, JC1106-11, JC1106-12, JC1106-13, JC1106-14, JC1106-15, JC1106-18, JC1106-19

6.2.1



CAS No.	Compound	Spike ug/l	BSP ug/l	BSP %	Limits
75-34-3	1,1-Dichloroethane	50	59.5	119	80-125
107-06-2	1,2-Dichloroethane	50	56.0	112	78-131
75-35-4	1,1-Dichloroethene	50	55.7	111	73-127
156-59-2	cis-1,2-Dichloroethene	50	52.4	105	77-118
156-60-5	trans-1,2-Dichloroethene	50	55.9	112	75-118
100-41-4	Ethylbenzene	50	52.3	105	80-118
75-09-2	Methylene chloride	50	57.0	114	75-122
127-18-4	Tetrachloroethene	50	50.8	102	69-138
108-88-3	Toluene	50	56.9	114	80-122
71-55-6	1,1,1-Trichloroethane	50	60.4	121	80-131
79-00-5	1,1,2-Trichloroethane	50	55.6	111	78-122
79-01-6	Trichloroethene	50	54.9	110	83-122
75-01-4	Vinyl chloride	50	53.8	108	57-138

CAS No.	Surrogate Recoveries	BSP	Limits
1868-53-7	Dibromofluoromethane	110%	76-120%
17060-07-0	1,2-Dichloroethane-D4	108%	73-122%
2037-26-5	Toluene-D8	103%	84-119%
460-00-4	4-Bromofluorobenzene	99%	78-117%

* = Outside of Control Limits.

Blank Spike Summary**Job Number:**

JC1106

Account:

UTC United Technologies Corporation

Project:

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3D4823-BS	3D112545.D	1	08/12/15	XC	n/a	n/a	V3D4823

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC1106-17

6.2.2



CAS No.	Compound	Spike ug/l	BSP ug/l	BSP %	Limits
75-34-3	1,1-Dichloroethane	50	56.5	113	80-125
107-06-2	1,2-Dichloroethane	50	55.6	111	78-131
75-35-4	1,1-Dichloroethene	50	53.0	106	73-127
156-59-2	cis-1,2-Dichloroethene	50	49.0	98	77-118
156-60-5	trans-1,2-Dichloroethene	50	53.1	106	75-118
100-41-4	Ethylbenzene	50	50.6	101	80-118
75-09-2	Methylene chloride	50	53.5	107	75-122
127-18-4	Tetrachloroethene	50	52.7	105	69-138
108-88-3	Toluene	50	55.7	111	80-122
71-55-6	1,1,1-Trichloroethane	50	58.1	116	80-131
79-00-5	1,1,2-Trichloroethane	50	54.3	109	78-122
79-01-6	Trichloroethene	50	53.2	106	83-122
75-01-4	Vinyl chloride	50	60.2	120	57-138

CAS No.	Surrogate Recoveries	BSP	Limits
1868-53-7	Dibromofluoromethane	108%	76-120%
17060-07-0	1,2-Dichloroethane-D4	105%	73-122%
2037-26-5	Toluene-D8	104%	84-119%
460-00-4	4-Bromofluorobenzene	102%	78-117%

* = Outside of Control Limits.

Blank Spike Summary

Job Number: JC1106
Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V3D4824-BS	3D112570.D	1	08/12/15	XC	n/a	n/a	V3D4824

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC1106-1, JC1106-2, JC1106-9, JC1106-16

6.2.3



CAS No.	Compound	Spike ug/l	BSP ug/l	BSP %	Limits
75-34-3	1,1-Dichloroethane	50	55.3	111	80-125
107-06-2	1,2-Dichloroethane	50	53.6	107	78-131
75-35-4	1,1-Dichloroethene	50	52.5	105	73-127
156-59-2	cis-1,2-Dichloroethene	50	49.1	98	77-118
156-60-5	trans-1,2-Dichloroethene	50	52.4	105	75-118
100-41-4	Ethylbenzene	50	49.3	99	80-118
75-09-2	Methylene chloride	50	53.2	106	75-122
127-18-4	Tetrachloroethene	50	48.7	97	69-138
108-88-3	Toluene	50	53.9	108	80-122
71-55-6	1,1,1-Trichloroethane	50	56.1	112	80-131
79-00-5	1,1,2-Trichloroethane	50	51.9	104	78-122
79-01-6	Trichloroethene	50	51.9	104	83-122
75-01-4	Vinyl chloride	50	55.2	110	57-138

CAS No.	Surrogate Recoveries	BSP	Limits
1868-53-7	Dibromofluoromethane	110%	76-120%
17060-07-0	1,2-Dichloroethane-D4	107%	73-122%
2037-26-5	Toluene-D8	103%	84-119%
460-00-4	4-Bromofluorobenzene	101%	78-117%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary**Job Number:** JC1106**Account:** UTC United Technologies Corporation**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
JC1106-13MS	3D112526.D	1	08/11/15	XC	n/a	n/a	V3D4822
JC1106-13MSD	3D112527.D	1	08/11/15	XC	n/a	n/a	V3D4822
JC1106-13	3D112524.D	1	08/11/15	XC	n/a	n/a	V3D4822

The QC reported here applies to the following samples:**Method:** SW846 8260C

JC1106-3, JC1106-4, JC1106-5, JC1106-6, JC1106-7, JC1106-8, JC1106-10, JC1106-11, JC1106-12, JC1106-13, JC1106-14, JC1106-15, JC1106-18, JC1106-19

CAS No.	Compound	JC1106-13		MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
		ug/l	Q							
75-34-3	1,1-Dichloroethane	1.6	50	60.8	118	50	57.4	112	6	60-129/13
107-06-2	1,2-Dichloroethane	ND	50	59.2	118	50	58.2	116	2	72-133/12
75-35-4	1,1-Dichloroethene	ND	50	54.2	108	50	50.2	100	8	40-137/17
156-59-2	cis-1,2-Dichloroethene	0.57	J	55.5	110	50	52.6	104	5	57-128/13
156-60-5	trans-1,2-Dichloroethene	ND	50	57.0	114	50	54.1	108	5	53-128/15
100-41-4	Ethylbenzene	ND	50	53.8	108	50	52.4	105	3	38-139/12
75-09-2	Methylene chloride	ND	50	59.3	119	50	56.5	113	5	63-128/13
127-18-4	Tetrachloroethene	0.43	J	49.2	98	50	48.3	96	2	43-145/15
108-88-3	Toluene	ND	50	58.1	116	50	55.9	112	4	51-136/13
71-55-6	1,1,1-Trichloroethane	3.0	50	59.4	113	50	56.1	106	6	51-141/16
79-00-5	1,1,2-Trichloroethane	ND	50	59.0	118	50	56.4	113	5	71-127/12
79-01-6	Trichloroethene	0.40	J	55.2	110	50	53.5	106	3	55-136/14
75-01-4	Vinyl chloride	ND	50	43.6	87	50	42.6	85	2	34-147/17

CAS No.	Surrogate Recoveries	MS	MSD	JC1106-13	Limits
1868-53-7	Dibromofluoromethane	110%	108%	111%	76-120%
17060-07-0	1,2-Dichloroethane-D4	107%	104%	112%	73-122%
2037-26-5	Toluene-D8	103%	103%	98%	84-119%
460-00-4	4-Bromofluorobenzene	100%	102%	97%	78-117%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
JC1072-1MS	3D112566.D	1	08/12/15	XC	n/a	n/a	V3D4823
JC1072-1MSD	3D112567.D	1	08/12/15	XC	n/a	n/a	V3D4823
JC1072-1	3D112560.D	1	08/12/15	XC	n/a	n/a	V3D4823

The QC reported here applies to the following samples:

Method: SW846 8260C

JC1106-17

CAS No.	Compound	JC1072-1 ug/l	Q	Spike ug/l	MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
75-34-3	1,1-Dichloroethane	ND		50	55.1	110	50	55.7	111	1	60-129/13
107-06-2	1,2-Dichloroethane	ND		50	56.2	112	50	58.0	116	3	72-133/12
75-35-4	1,1-Dichloroethene	ND		50	50.6	101	50	50.6	101	0	40-137/17
156-59-2	cis-1,2-Dichloroethene	ND		50	49.7	99	50	50.5	101	2	57-128/13
156-60-5	trans-1,2-Dichloroethene	ND		50	53.0	106	50	53.2	106	0	53-128/15
100-41-4	Ethylbenzene	ND		50	49.3	99	50	51.3	103	4	38-139/12
75-09-2	Methylene chloride	ND		50	54.7	109	50	54.7	109	0	63-128/13
127-18-4	Tetrachloroethene	ND		50	46.1	92	50	48.5	97	5	43-145/15
108-88-3	Toluene	ND		50	55.3	111	50	56.0	112	1	51-136/13
71-55-6	1,1,1-Trichloroethane	ND		50	55.2	110	50	54.4	109	1	51-141/16
79-00-5	1,1,2-Trichloroethane	ND		50	55.7	111	50	58.2	116	4	71-127/12
79-01-6	Trichloroethene	ND		50	52.6	105	50	53.8	108	2	55-136/14
75-01-4	Vinyl chloride	ND		50	54.9	110	50	56.2	112	2	34-147/17

CAS No.	Surrogate Recoveries	MS	MSD	JC1072-1	Limits
1868-53-7	Dibromofluoromethane	110%	107%	108%	76-120%
17060-07-0	1,2-Dichloroethane-D4	106%	105%	112%	73-122%
2037-26-5	Toluene-D8	105%	103%	100%	84-119%
460-00-4	4-Bromofluorobenzene	100%	99%	95%	78-117%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
JC1106-9MS	3D112580.D	1	08/12/15	XC	n/a	n/a	V3D4824
JC1106-9MSD	3D112581.D	1	08/12/15	XC	n/a	n/a	V3D4824
JC1106-9	3D112572.D	1	08/12/15	XC	n/a	n/a	V3D4824

The QC reported here applies to the following samples:

Method: SW846 8260C

JC1106-1, JC1106-2, JC1106-9, JC1106-16

CAS No.	Compound	JC1106-9 ug/l	Spike Q	MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits Rec/RPD
75-34-3	1,1-Dichloroethane	1.1		53.7	105	50	52.5	103	2	60-129/13
107-06-2	1,2-Dichloroethane	ND		54.0	108	50	53.1	106	2	72-133/12
75-35-4	1,1-Dichloroethene	0.93	J	50	49.7	98	50	48.8	96	2
156-59-2	cis-1,2-Dichloroethene	3.5		54.0	101	50	52.5	98	3	57-128/13
156-60-5	trans-1,2-Dichloroethene	ND		51.0	102	50	50.0	100	2	53-128/15
100-41-4	Ethylbenzene	ND		50.9	102	50	50.5	101	1	38-139/12
75-09-2	Methylene chloride	ND		54.1	108	50	53.2	106	2	63-128/13
127-18-4	Tetrachloroethene	ND		47.0	94	50	47.5	95	1	43-145/15
108-88-3	Toluene	ND		53.6	107	50	54.2	108	1	51-136/13
71-55-6	1,1,1-Trichloroethane	53.6		78.6	50* a	50	78.5	50* a	0	51-141/16
79-00-5	1,1,2-Trichloroethane	ND		54.7	109	50	55.0	110	1	71-127/12
79-01-6	Trichloroethene	2.6		52.3	99	50	52.2	99	0	55-136/14
75-01-4	Vinyl chloride	ND		49.1	98	50	48.9	98	0	34-147/17

CAS No.	Surrogate Recoveries	MS	MSD	JC1106-9	Limits
1868-53-7	Dibromofluoromethane	107%	104%	112%	76-120%
17060-07-0	1,2-Dichloroethane-D4	98%	97%	111%	73-122%
2037-26-5	Toluene-D8	102%	102%	102%	84-119%
460-00-4	4-Bromofluorobenzene	103%	101%	96%	78-117%

(a) Outside control limits due to matrix interference.

* = Outside of Control Limits.

Instrument Performance Check (BFB)

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample: V3D4810-BFB

Injection Date: 07/29/15

Lab File ID: 3D112234.D

Injection Time: 19:16

Instrument ID: GCMS3D

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	15.0 - 40.0% of mass 95	9449	16.9	Pass
75	30.0 - 60.0% of mass 95	24824	44.4	Pass
95	Base peak, 100% relative abundance	55917	100.0	Pass
96	5.0 - 9.0% of mass 95	3622	6.48	Pass
173	Less than 2.0% of mass 174	0	0.00	(0.00) ^a Pass
174	50.0 - 120.0% of mass 95	55002	98.4	Pass
175	5.0 - 9.0% of mass 174	4400	7.87	(8.00) ^a Pass
176	95.0 - 101.0% of mass 174	53496	95.7	(97.3) ^a Pass
177	5.0 - 9.0% of mass 176	3680	6.58	(6.88) ^b Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V3D4810-IC4810	3D112235.D	07/29/15	20:15	00:59	Initial cal 0.2
V3D4810-IC4810	3D112236.D	07/29/15	20:42	01:26	Initial cal 0.5
V3D4810-IC4810	3D112237.D	07/29/15	21:10	01:54	Initial cal 1
V3D4810-IC4810	3D112238.D	07/29/15	21:37	02:21	Initial cal 2
V3D4810-IC4810	3D112239.D	07/29/15	22:04	02:48	Initial cal 5
V3D4810-IC4810	3D112240.D	07/29/15	22:31	03:15	Initial cal 10
V3D4810-IC4810	3D112241.D	07/29/15	22:58	03:42	Initial cal 20
V3D4810-ICC4810	3D112242.D	07/29/15	23:25	04:09	Initial cal 50
V3D4810-IC4810	3D112243.D	07/29/15	23:53	04:37	Initial cal 100
V3D4810-IC4810	3D112244.D	07/30/15	00:20	05:04	Initial cal 200
V3D4810-ICV4810	3D112247.D	07/30/15	01:41	06:25	Initial cal verification 50

Instrument Performance Check (BFB)**Job Number:**

JC1106

Account:

UTC United Technologies Corporation

Project:

ENSRLW: UTAS Plants 1/2 Facility, Rockford, IL

Sample: V3D4822-BFB**Injection Date:** 08/11/15**Lab File ID:** 3D112515.D**Injection Time:** 09:53**Instrument ID:** GCMS3D

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	15.0 - 40.0% of mass 95	8665	17.8	Pass
75	30.0 - 60.0% of mass 95	21568	44.2	Pass
95	Base peak, 100% relative abundance	48757	100.0	Pass
96	5.0 - 9.0% of mass 95	3479	7.14	Pass
173	Less than 2.0% of mass 174	0	0.00	(0.00) ^a Pass
174	50.0 - 120.0% of mass 95	49194	100.9	Pass
175	5.0 - 9.0% of mass 174	3521	7.22	(7.16) ^a Pass
176	95.0 - 101.0% of mass 174	47664	97.8	(96.9) ^a Pass
177	5.0 - 9.0% of mass 176	3271	6.71	(6.86) ^b Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V3D4822-CC4810	3D112517.D	08/11/15	11:16	01:23	Continuing cal 20
V3D4822-MB	3D112519.D	08/11/15	12:38	02:45	Method Blank
V3D4822-BS	3D112520.D	08/11/15	13:12	03:19	Blank Spike
JC1106-7	3D112522.D	08/11/15	14:09	04:16	HSSER-FBLK01-080415
JC1106-5	3D112523.D	08/11/15	14:36	04:43	HSSER-MW203-080415
JC1106-13	3D112524.D	08/11/15	15:03	05:10	HSSER-GMZ02-080515
JC1106-10	3D112525.D	08/11/15	15:30	05:37	HSSER-SMW21-080515
JC1106-13MS	3D112526.D	08/11/15	16:03	06:10	Matrix Spike
JC1106-13MSD	3D112527.D	08/11/15	16:30	06:37	Matrix Spike Duplicate
JC1106-19	3D112529.D	08/11/15	17:24	07:31	HSSER-TRIP01-080415
JC1106-6	3D112530.D	08/11/15	17:51	07:58	HSSER-MW07FGA-080515
JC1106-3	3D112531.D	08/11/15	18:18	08:25	HSSER-SMW01-080415
JC1106-4	3D112532.D	08/11/15	18:45	08:52	HSSER-SMW02-080415
JC1106-15	3D112533.D	08/11/15	19:13	09:20	HSSER-PMW01-080615
JC1106-18	3D112534.D	08/11/15	19:40	09:47	HSSER-SMW04-080615
JC1106-8	3D112535.D	08/11/15	20:07	10:14	HSSER-SMW19-080515
JC1106-14	3D112536.D	08/11/15	20:34	10:41	HSSER-DUP01-080515
JC1106-12	3D112537.D	08/11/15	21:01	11:08	HSSER-GMZ03-080515
JC1106-11	3D112538.D	08/11/15	21:28	11:35	HSSER-SMW20-080515

6.4.2



Instrument Performance Check (BFB)

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample: V3D4823-BFB

Injection Date: 08/11/15

Lab File ID: 3D112541.D

Injection Time: 22:50

Instrument ID: GCMS3D

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	15.0 - 40.0% of mass 95	7142	16.6	Pass
75	30.0 - 60.0% of mass 95	19101	44.5	Pass
95	Base peak, 100% relative abundance	42928	100.0	Pass
96	5.0 - 9.0% of mass 95	2834	6.60	Pass
173	Less than 2.0% of mass 174	632	1.47	(1.36) ^a
174	50.0 - 120.0% of mass 95	46456	108.2	Pass
175	5.0 - 9.0% of mass 174	3302	7.69	(7.11) ^a
176	95.0 - 101.0% of mass 174	44520	103.7	(95.8) ^a
177	5.0 - 9.0% of mass 176	2956	6.89	(6.64) ^b

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V3D4823-CC4810	3D112542.D	08/11/15	23:17	00:27	Continuing cal 50
V3D4823-MB	3D112544.D	08/12/15	00:11	01:21	Method Blank
V3D4823-BS	3D112545.D	08/12/15	00:38	01:48	Blank Spike
ZZZZZZ	3D112549.D	08/12/15	02:26	03:36	(unrelated sample)
ZZZZZZ	3D112550.D	08/12/15	02:53	04:03	(unrelated sample)
ZZZZZZ	3D112551.D	08/12/15	03:20	04:30	(unrelated sample)
ZZZZZZ	3D112552.D	08/12/15	03:47	04:57	(unrelated sample)
ZZZZZZ	3D112553.D	08/12/15	04:13	05:23	(unrelated sample)
JC1106-17	3D112555.D	08/12/15	05:07	06:17	HSSER-EBLK01-080615
JC1072-1	3D112560.D	08/12/15	07:22	08:32	(used for QC only; not part of job JC1106)
ZZZZZZ	3D112561.D	08/12/15	07:49	08:59	(unrelated sample)

Instrument Performance Check (BFB)

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample: V3D4824-BFB

Injection Date: 08/12/15

Lab File ID: 3D112562.D

Injection Time: 08:16

Instrument ID: GCMS3D

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	15.0 - 40.0% of mass 95	8422	18.2	Pass
75	30.0 - 60.0% of mass 95	20373	44.1	Pass
95	Base peak, 100% relative abundance	46160	100.0	Pass
96	5.0 - 9.0% of mass 95	2967	6.43	Pass
173	Less than 2.0% of mass 174	385	0.83	(0.81) ^a Pass
174	50.0 - 120.0% of mass 95	47440	102.8	Pass
175	5.0 - 9.0% of mass 174	3561	7.71	(7.51) ^a Pass
176	95.0 - 101.0% of mass 174	46184	100.1	(97.4) ^a Pass
177	5.0 - 9.0% of mass 176	2874	6.23	(6.22) ^b Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V3D4824-CC4810	3D112563.D	08/12/15	09:17	01:01	Continuing cal 20
V3D4823-MB2	3D112565.D	08/12/15	10:15	01:59	Method Blank
JC1072-1MS	3D112566.D	08/12/15	10:44	02:28	Matrix Spike
JC1072-1MSD	3D112567.D	08/12/15	11:11	02:55	Matrix Spike Duplicate
V3D4824-MB	3D112569.D	08/12/15	12:06	03:50	Method Blank
V3D4824-BS	3D112570.D	08/12/15	12:33	04:17	Blank Spike
JC1106-9	3D112572.D	08/12/15	13:27	05:11	HSSER-GMZ04-080515
JC1106-16	3D112573.D	08/12/15	13:54	05:38	HSSER-PMW02-080615
JC1106-2	3D112574.D	08/12/15	14:21	06:05	HSSER-GMZ01-080415
JC1106-1	3D112575.D	08/12/15	14:48	06:32	HSSER-SMW08-080415
ZZZZZZ	3D112576.D	08/12/15	15:15	06:59	(unrelated sample)
ZZZZZZ	3D112577.D	08/12/15	15:42	07:26	(unrelated sample)
ZZZZZZ	3D112578.D	08/12/15	16:10	07:54	(unrelated sample)
ZZZZZZ	3D112579.D	08/12/15	16:37	08:21	(unrelated sample)
JC1106-9MS	3D112580.D	08/12/15	17:04	08:48	Matrix Spike
JC1106-9MSD	3D112581.D	08/12/15	17:31	09:15	Matrix Spike Duplicate
ZZZZZZ	3D112582.D	08/12/15	17:58	09:42	(unrelated sample)
ZZZZZZ	3D112583.D	08/12/15	18:25	10:09	(unrelated sample)
ZZZZZZ	3D112584.D	08/12/15	18:52	10:36	(unrelated sample)

Volatile Internal Standard Area Summary

Page 1 of 1

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Check Std:	V3D4822-CC4810	Injection Date:	08/11/15
Lab File ID:	3D112517.D	Injection Time:	11:16
Instrument ID:	GCMS3D	Method:	SW846 8260C

	IS 1 AREA	IS 2 AREA	IS 3 AREA	IS 4 AREA	IS 5 AREA	RT	IS 1 AREA	IS 2 AREA	IS 3 AREA	IS 4 AREA	IS 5 AREA	RT
Check Std	105534	7.16	204849	9.39	234799	10.31	194668	13.48	122430	15.79		
Upper Limit ^a	211068	7.66	409698	9.89	469598	10.81	389336	13.98	244860	16.29		
Lower Limit ^b	52767	6.66	102425	8.89	117400	9.81	97334	12.98	61215	15.29		

Lab Sample ID	IS 1 AREA	IS 2 AREA	IS 3 AREA	IS 4 AREA	IS 5 AREA	RT	IS 1 AREA	IS 2 AREA	IS 3 AREA	IS 4 AREA	IS 5 AREA	RT
V3D4822-MB	103952	7.18	226869	9.39	264986	10.31	216656	13.48	129511	15.79		
V3D4822-BS	107290	7.18	202270	9.39	233933	10.31	196021	13.48	124533	15.79		
JC1106-7	92150	7.17	192784	9.39	222789	10.31	184238	13.48	117577	15.79		
JC1106-5	97975	7.17	195078	9.39	228938	10.31	186093	13.48	118621	15.79		
JC1106-13	91266	7.17	186566	9.40	219744	10.31	177389	13.48	110292	15.79		
JC1106-10	87229	7.17	179050	9.39	210954	10.31	172979	13.48	107549	15.79		
JC1106-13MS	109061	7.17	200873	9.39	227802	10.31	191425	13.48	118628	15.79		
JC1106-13MSD	106157	7.18	209488	9.39	234394	10.31	194946	13.48	117427	15.79		
JC1106-19	97633	7.17	205375	9.40	241116	10.31	197855	13.48	123858	15.79		
JC1106-6	95232	7.17	188915	9.39	220873	10.31	182140	13.48	113662	15.79		
JC1106-3	93947	7.17	191420	9.39	223101	10.31	183270	13.48	111855	15.79		
JC1106-4	93695	7.17	187081	9.40	219575	10.31	184978	13.48	115133	15.79		
JC1106-15	94051	7.17	193794	9.39	228111	10.31	188105	13.48	116304	15.79		
JC1106-18	88719	7.17	186366	9.39	213690	10.31	175269	13.48	108187	15.79		
JC1106-8	99250	7.17	189230	9.39	220375	10.31	181092	13.48	111525	15.79		
JC1106-14	91107	7.17	183583	9.39	216133	10.31	177301	13.48	109805	15.79		
JC1106-12	84872	7.17	176303	9.39	208560	10.31	169036	13.48	106176	15.79		
JC1106-11	84171	7.17	181857	9.39	212162	10.31	174691	13.48	107285	15.79		

IS 1 = Tert Butyl Alcohol-D9

IS 2 = Pentafluorobenzene

IS 3 = 1,4-Difluorobenzene

IS 4 = Chlorobenzene-D5

IS 5 = 1,4-Dichlorobenzene-d4

(a) Upper Limit = + 100% of check standard area; Retention time + 0.5 minutes.

(b) Lower Limit = -50% of check standard area; Retention time -0.5 minutes.

Volatile Internal Standard Area Summary

Page 1 of 1

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Check Std:	V3D4823-CC4810	Injection Date:	08/11/15
Lab File ID:	3D112542.D	Injection Time:	23:17
Instrument ID:	GCMS3D	Method:	SW846 8260C

	IS 1 AREA	IS 2 RT	IS 2 AREA	IS 2 RT	IS 3 AREA	IS 3 RT	IS 4 AREA	IS 4 RT	IS 5 AREA	IS 5 RT
Check Std	90136	7.17	190390	9.39	217047	10.31	183008	13.48	113415	15.79
Upper Limit ^a	180272	7.67	380780	9.89	434094	10.81	366016	13.98	226830	16.29
Lower Limit ^b	45068	6.67	95195	8.89	108524	9.81	91504	12.98	56708	15.29

Lab Sample ID	IS 1 AREA	IS 1 RT	IS 2 AREA	IS 2 RT	IS 3 AREA	IS 3 RT	IS 4 AREA	IS 4 RT	IS 5 AREA	IS 5 RT
V3D4823-MB	82447	7.16	171750	9.39	198901	10.31	165044	13.48	101680	15.79
V3D4823-BS	93840	7.17	196820	9.40	222078	10.31	188570	13.48	114311	15.79
ZZZZZZ	95247	7.17	198301	9.39	228803	10.31	187554	13.48	115293	15.79
ZZZZZZ	93410	7.16	194718	9.39	227275	10.31	186257	13.48	117632	15.79
ZZZZZZ	93613	7.16	194392	9.39	224971	10.31	186632	13.48	115937	15.79
ZZZZZZ	91157	7.17	191813	9.39	226390	10.31	185266	13.48	113424	15.79
ZZZZZZ	89555	7.17	186162	9.39	217020	10.31	179683	13.48	114175	15.79
JC1106-17	94319	7.17	190266	9.39	224022	10.31	184821	13.48	114314	15.79
JC1072-1	90352	7.16	187714	9.39	220418	10.31	181445	13.48	111052	15.79
ZZZZZZ	88952	7.17	180001	9.39	212568	10.31	174775	13.48	108634	15.79

IS 1 = Tert Butyl Alcohol-D9

IS 2 = Pentafluorobenzene

IS 3 = 1,4-Difluorobenzene

IS 4 = Chlorobenzene-D5

IS 5 = 1,4-Dichlorobenzene-d4

(a) Upper Limit = + 100% of check standard area; Retention time + 0.5 minutes.

(b) Lower Limit = -50% of check standard area; Retention time -0.5 minutes.

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Volatile Internal Standard Area Summary

Page 1 of 1

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Check Std:	V3D4824-CC4810	Injection Date:	08/12/15
Lab File ID:	3D112563.D	Injection Time:	09:17
Instrument ID:	GCMS3D	Method:	SW846 8260C

	IS 1 AREA	RT	IS 2 AREA	RT	IS 3 AREA	RT	IS 4 AREA	RT	IS 5 AREA	RT
Check Std	97817	7.17	204794	9.40	227986	10.32	187452	13.48	110302	15.79
Upper Limit ^a	195634	7.67	409588	9.90	455972	10.82	374904	13.98	220604	16.29
Lower Limit ^b	48909	6.67	102397	8.90	113993	9.82	93726	12.98	55151	15.29

Lab Sample ID	IS 1 AREA	RT	IS 2 AREA	RT	IS 3 AREA	RT	IS 4 AREA	RT	IS 5 AREA	RT
V3D4823-MB2	89468	7.18	182832	9.39	209972	10.31	174088	13.48	109833	15.79
JC1072-1MS	99162	7.18	201673	9.39	227956	10.31	197596	13.48	122225	15.79
JC1072-1MSD	99080	7.17	197615	9.39	223819	10.31	186124	13.48	115672	15.79
V3D4824-MB	95029	7.18	192060	9.39	221163	10.31	183575	13.48	114634	15.79
V3D4824-BS	97953	7.17	207104	9.39	238560	10.31	200201	13.48	122169	15.79
JC1106-9	91436	7.19	193182	9.39	221437	10.31	184042	13.48	114924	15.79
JC1106-16	88764	7.18	186964	9.39	217383	10.31	181788	13.48	113725	15.79
JC1106-2	92619	7.17	180692	9.39	212538	10.31	175332	13.48	108820	15.79
JC1106-1	93640	7.18	196615	9.39	224320	10.31	180779	13.48	111990	15.79
ZZZZZZ	103650	7.17	196401	9.39	225527	10.31	186617	13.48	114903	15.79
ZZZZZZ	107002	7.17	204398	9.39	232631	10.31	188859	13.48	114136	15.79
ZZZZZZ	154553	7.16	223480	9.40	250790	10.31	206024	13.48	124766	15.79
ZZZZZZ	145130	7.16	230897	9.39	261778	10.31	212455	13.48	132696	15.79
JC1106-9MS	118942	7.18	243673	9.40	269432	10.31	220893	13.48	129859	15.79
JC1106-9MSD	120440	7.18	245404	9.39	271506	10.31	223373	13.48	132733	15.79
ZZZZZZ	113579	7.18	239484	9.39	274644	10.31	224882	13.48	143888	15.79
ZZZZZZ	114948	7.17	230024	9.39	265060	10.31	216179	13.48	132533	15.79
ZZZZZZ	163458	7.16	211232	9.39	244674	10.31	197571	13.48	122611	15.79

IS 1 = Tert Butyl Alcohol-D9

IS 2 = Pentafluorobenzene

IS 3 = 1,4-Difluorobenzene

IS 4 = Chlorobenzene-D5

IS 5 = 1,4-Dichlorobenzene-d4

(a) Upper Limit = + 100% of check standard area; Retention time + 0.5 minutes.

(b) Lower Limit = -50% of check standard area; Retention time -0.5 minutes.

Volatile Surrogate Recovery Summary

Page 1 of 1

Job Number: JC1106

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Method: SW846 8260C

Matrix: AQ

Samples and QC shown here apply to the above method

Lab Sample ID	Lab File ID	S1	S2	S3	S4
JC1106-1	3D112575.D	112	110	99	97
JC1106-2	3D112574.D	112	115	100	95
JC1106-3	3D112531.D	110	113	101	98
JC1106-4	3D112532.D	114	113	101	97
JC1106-5	3D112523.D	112	114	101	97
JC1106-6	3D112530.D	111	112	100	96
JC1106-7	3D112522.D	111	112	101	96
JC1106-8	3D112535.D	113	115	101	97
JC1106-9	3D112572.D	112	111	102	96
JC1106-10	3D112525.D	113	114	99	97
JC1106-11	3D112538.D	114	113	101	97
JC1106-12	3D112537.D	113	115	99	97
JC1106-13	3D112524.D	111	112	98	97
JC1106-14	3D112536.D	114	117	101	97
JC1106-15	3D112533.D	113	112	100	96
JC1106-16	3D112573.D	111	112	100	95
JC1106-17	3D112555.D	108	112	101	98
JC1106-18	3D112534.D	111	112	100	98
JC1106-19	3D112529.D	112	113	101	96
JC1072-1MS	3D112566.D	110	106	105	100
JC1072-1MSD	3D112567.D	107	105	103	99
JC1106-13MS	3D112526.D	110	107	103	100
JC1106-13MSD	3D112527.D	108	104	103	102
JC1106-9MS	3D112580.D	107	98	102	103
JC1106-9MSD	3D112581.D	104	97	102	101
V3D4822-BS	3D112520.D	110	108	103	99
V3D4822-MB	3D112519.D	108	112	102	100
V3D4823-BS	3D112545.D	108	105	104	102
V3D4823-MB	3D112544.D	112	114	102	96
V3D4824-BS	3D112570.D	110	107	103	101
V3D4824-MB	3D112569.D	114	110	101	97
V3D4823-MB2	3D112565.D	113	112	102	95

Surrogate Compounds

Recovery Limits

S1 = Dibromofluoromethane	76-120%
S2 = 1,2-Dichloroethane-D4	73-122%
S3 = Toluene-D8	84-119%
S4 = 4-Bromofluorobenzene	78-117%

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6
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Initial Calibration Summary

Page 1 of 5

Job Number: JC1106

Sample: V3D4810-ICC4810
 Lab FileID: 3D112242.D

Account: UTC United Technologies Corporation
 Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Response Factor Report MS3D

Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)
 Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 Last Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Calibration Files

5 =3D112239.D	0.5 =3D112236.D	2 =3D112238.D	50 =3D112242.D
100 =3D112243.D	1 =3D112237.D	200 =3D112244.D	20 =3D112241.D
10 =3D112240.D	0.2 =3D112235.D	=	=

Compound

	5	0.5	2	50	100	1	200	20	10	0.2	Avg	%RSD
--	---	-----	---	----	-----	---	-----	----	----	-----	-----	------

1) I Tert Butyl Alcohol-d9	-----ISTD-----										
2) 1,4-dioxane	0.077 0.076 0.089 0.088 0.066 0.084 0.088 0.079										
3) tertiary butyl alcohol	1.062 1.166 1.077 1.068 1.038 1.109 1.037										
4) I pentafluorobenzene	-----ISTD-----										
5) FREON 115	0.000# -1.00										
6) FREON 152A	0.000# -1.00										
7) chlorotrifluoroethene	0.000# -1.00										
8) chlorodifluoromethane	0.683 0.700 0.677 0.697 0.696 0.686 0.613										
9) dichlorodifluoromethane	0.983 0.922 0.904 0.888 0.816 0.859 0.946 0.953										
10) chloromethane	0.687 0.843 0.635 0.679 0.688 0.671 0.734 0.649 0.646 0.993 0.723										
11) vinyl chloride	15.54 0.741 0.758 0.680 0.755 0.784 0.687 0.830 0.716 0.724 0.572 0.725										
12) bromomethane	9.63 0.472 0.564 0.447 0.449 0.441 0.501 0.422 0.451 0.447 0.489 0.468										
13) chloroethane	8.78 0.319 0.392 0.290 0.293 0.283 0.321 0.276 0.288 0.301										
14) 1,3-butadiene	11.55 0.000# -1.00										
15) vinyl bromide	0.000# -1.00										
16) trichlorofluoromethane	5.56 0.808 0.708 0.737 0.786 0.774 0.674 0.768 0.778 0.772										
17) pentane	0.000# -1.00										
18) ethyl ether	8.81 0.179 0.180 0.178 0.179 0.227 0.181 0.187 0.182										
19) freon 123a	0.000# -1.00										
20) FREON 123	0.000# -1.00										
21) acrolein	4.73 0.080 0.077 0.075 0.068 0.077 0.075 0.074 0.074										
22) 1,1-dichloroethene	12.98 0.587 0.419 0.602 0.620 0.600 0.720 0.605 0.609 0.593										
23) acetone	0.595 0.595 12.98										

Initial Calibration Summary

Page 2 of 5

Job Number: JC1106

Sample: V3D4810-ICC4810
Lab FileID: 3D112242.DAccount: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	0.120	0.095	0.109	0.109	0.120	0.120	0.112	9.00					
24)	allyl chloride	0.182	0.127	0.180	0.201	0.190	0.202	0.191	0.203	0.187	0.185	12.62	
25)	acetonitrile	0.038	0.037	0.035	0.038	0.041	0.038	0.037	0.034	0.037	5.70		
26)	acetalddehyde									0.000#	-1.00		
27)	iodomethane	0.836	0.744	0.873	0.911	0.891	0.986	0.905	0.904	0.857	0.550	0.846	14.32
28)	iso-butyl alcohol	0.013	0.012	0.012	0.013		0.014	0.012	0.012		0.012	6.76	
29)	carbon disulfide	1.483	1.330	1.542	1.564	1.521	1.794	1.534	1.569	1.489	1.213	1.504	10.16
30)	methylene chloride	0.425	0.452	0.449	0.454	0.440	0.517	0.441	0.456	0.440	0.382	0.446	7.46
31)	methyl acetate	0.279	0.250	0.252	0.256	0.194	0.270	0.252	0.237		0.249	10.25	
32)	methyl tert butyl ether	1.250	1.300	1.272	1.319	1.333	1.469	1.343	1.370	1.315	1.229	1.320	5.13
33)	trans-1,2-dichloroethene	0.524	0.507	0.570	0.561	0.542	0.656	0.544	0.569	0.531	0.423	0.543	10.74
34)	di-isopropyl ether	1.350	1.555	1.330	1.366	1.365	1.560	1.375	1.343	1.229	1.502	1.396	7.44
35)	ethyl tert-butyl ether	1.376	1.366	1.278	1.376	1.404	1.397	1.427	1.344	1.209	0.972	1.315	10.41
36)	2-butanone	0.026	0.034	0.039		0.040	0.039	0.035		0.036		14.42	
37)	1,1-dichloroethane	0.636	0.569	0.637	0.682	0.667	0.757	0.668	0.696	0.658	0.477	0.645	11.76
38)	chloroprene	0.487	0.445	0.515	0.494	0.485	0.558	0.488	0.465	0.426	0.355	0.472	11.58
39)	acrylonitrile	0.125	0.124	0.117	0.133	0.139	0.141	0.142	0.144	0.137	0.089	0.129	12.85
40)	vinyl acetate	0.039	0.050	0.054		0.056	0.050	0.038		0.048		15.71	
41)	ethyl acetate	0.054	0.061	0.043	0.045		0.045	0.051	0.044		0.049		13.52
42)	2,2-dichloropropane	0.625	0.710	0.656	0.643	0.643	0.803	0.634	0.658	0.634	0.739	0.675	8.61
43)	cis-1,2-dichloroethene	0.409	0.497	0.462	0.416	0.417	0.533	0.423	0.430	0.417	0.354	0.436	11.47
44)	propionitrile	0.045	0.044	0.047	0.051	0.044	0.054	0.051	0.048		0.048		7.90
45)	methyl acrylate	0.284	0.249	0.329	0.350		0.368	0.362	0.326		0.324		13.39
46)	bromochloromethane	0.198	0.159	0.194	0.217	0.220	0.238	0.222	0.227	0.212		0.210	11.12
47)	tetrahydrofuran	0.120	0.156	0.113	0.116		0.124	0.130	0.125		0.126		11.22
48)	chloroform	0.586	0.616	0.636	0.638	0.630	0.695	0.641	0.638	0.595	0.543	0.622	6.51
49)	Tert-Butyl Formate	0.352	0.379	0.333	0.372	0.389	0.345	0.405	0.363	0.318		0.362	7.66
50)	dibromofluoromethane (s)	0.379	0.385	0.377	0.386	0.385	0.386	0.397	0.381	0.378	0.376	0.383	1.65
51)	1,2-dichloroethane-d4 (s)	0.361	0.361	0.357	0.350	0.348	0.357	0.357	0.353	0.350	0.356	0.355	1.29
52)	freon 113	0.383	0.367	0.374	0.362	0.396	0.349	0.364	0.334		0.366		5.24
53)	methacrylonitrile												

Initial Calibration Summary

Page 3 of 5

Job Number: JC1106

Sample: V3D4810-ICC4810
Lab FileID: 3D112242.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	0.162	0.206	0.181	0.195	0.206	0.193	0.172	0.188	9.03
54)	1,1,1-trichloroethane							0.642	6.43
	0.600	0.591	0.609	0.657	0.658	0.719	0.662	0.666	0.615
55)	tert amyl alcohol							0.000#	-1.00
56)	2,2,4-Trimethylpentane							0.000#	-1.00
57)	tert-amyl methyl ether							0.000#	-1.00
	1.290	1.548	1.277	1.279	1.318	1.375	1.360	1.259	1.130
								1.305	1.314
									8.05
58)	I 1,4-difluorobenzene	-----	-----	-----	-----	-----	-----		
59)	epichlorohydrin							0.027	8.27
	0.028	0.030	0.024	0.027	0.031	0.027	0.027	0.026	
60)	n-butyl alcohol							0.008#	10.11
	0.008	0.007	0.008	0.009	0.009	0.009	0.008	0.008	
61)	cyclohexane							0.586	6.93
	0.544	0.568	0.593	0.589	0.679	0.562	0.577	0.573	
62)	carbon tetrachloride							0.484	13.04
	0.477	0.355	0.476	0.530	0.530	0.560	0.511	0.518	
63)	1,1-dichloropropene							0.360	13.77
	0.355	0.236	0.346	0.386	0.387	0.396	0.385	0.389	
64)	hexane ** This compound does not meet Initial Calibration criteria							0.322	9.18
	0.338	0.313	0.360	0.322	0.308	0.381	0.293	0.304	
								0.290	
								0.306	
65)	benzene							0.284	6.34
	1.113	1.136	1.187	1.193	1.201	1.278	1.210	1.207	
								1.155	
								1.070	
								1.175	4.98
66)	heptane							0.183	5.52
	0.193	0.195	0.181	0.179	0.194	0.171	0.178	0.170	
67)	isopropyl acetate							0.511	8.82
	0.501	0.523	0.515	0.558	0.441	0.570	0.525	0.455	
68)	1,2-dichloroethane							0.351	7.37
	0.331	0.292	0.340	0.363	0.364	0.369	0.364	0.374	
69)	trichloroethene							0.284	6.34
	0.258	0.275	0.298	0.286	0.286	0.320	0.290	0.289	
70)	ethyl acrylate							0.000#	-1.00
71)	tert amyl ethyl ether							0.000#	-1.00
72)	2-nitropropane							0.104	6.98
	0.110	0.097	0.105		0.102	0.114	0.095		
73)	2-chloroethyl vinyl ether							0.180	4.24
	0.180	0.174	0.182	0.190	0.178	0.181	0.190	0.167	
74)	methyl methacrylate							0.402	12.17
	0.355	0.319	0.413	0.444		0.445	0.440	0.395	
75)	1,2-dichloropropane							0.307	5.24
	0.292	0.292	0.290	0.312	0.313	0.340	0.311	0.315	
76)	methylcyclohexane							0.531	17.07
	0.573	0.391	0.585	0.594	0.581	0.618	0.568	0.552	
								0.502	
								0.350	
77)	dibromomethane							0.374	7.80
	0.172	0.128	0.180	0.188	0.191	0.191	0.190	0.199	
								0.187	
78)	bromodichloromethane							0.181	11.72
	0.351	0.369	0.354	0.387	0.388	0.417	0.396	0.392	
								0.371	
79)	cis-1,3-dichloropropene							0.313	
	0.404	0.356	0.436	0.466	0.476	0.458	0.463	0.485	
								0.448	
80)	toluene-d8 (s)							0.393	9.43
	1.088	1.089	1.105	1.101	1.093	1.106	1.069	1.102	
								1.096	
81)	4-methyl-2-pentanone							1.102	1.03
	0.108	0.104	0.110	0.122	0.092	0.126	0.120	0.112	
82)	toluene							0.112	9.72
	0.620	0.538	0.645	0.694	0.695	0.730	0.681	0.716	
								0.655	
83)	3-methyl-1-butanol							0.483	12.31
								0.646	

Initial Calibration Summary

Page 4 of 5

Job Number: JC1106

Sample: V3D4810-ICC4810
Lab FileID: 3D112242.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

0.008	0.008	0.008	0.008	0.010	0.009	0.008	0.008	0.008	0.008#	7.78
84) trans-1,3-dichloropropene	0.372	0.351	0.362	0.416	0.429	0.429	0.414	0.431	0.403	0.293
85) ethyl methacrylate	0.313	0.298	0.351	0.373	0.350	0.364	0.366	0.343	0.345	7.62
86) 1,1,2-trichloroethane	0.199	0.191	0.207	0.217	0.222	0.256	0.217	0.229	0.214	0.217
87) 2-hexanone	0.086	0.071	0.088	0.103		0.104	0.100	0.088	0.091	12.66
88) I chlorobenzene-d5										ISTD-----
89) tetrachloroethene	0.388	0.371	0.415	0.441	0.432	0.478	0.436	0.460	0.411	0.426
90) 1,3-dichloropropane	0.450	0.423	0.494	0.495	0.503	0.495	0.503	0.517	0.485	0.485
91) butyl acetate	0.187	0.173	0.205	0.222		0.234	0.202	0.184	0.201	10.82
92) 3,3-Dimethyl-1-Butanol	0.045	0.051	0.048	0.040	0.049	0.059	0.059	0.042	0.040	0.048
93) dibromochloromethane	0.354	0.267	0.357	0.404	0.418	0.381	0.426	0.416	0.381	0.378
94) 1,2-dibromoethane	0.304	0.260	0.296	0.333	0.342	0.301	0.347	0.357	0.331	0.319
95) n-Butyl Ether	1.412	1.485	1.514	1.511	1.670	1.499	1.558	1.452	1.513	5.10
96) chlorobenzene	0.871	0.815	0.883	0.940	0.953	0.992	0.962	0.980	0.890	0.681
97) 1,1,1,2-tetrachloroethane	0.360	0.287	0.376	0.428	0.437	0.417	0.470	0.422	0.382	0.398
98) ethylbenzene	1.442	1.493	1.498	1.566	1.581	1.747	1.590	1.612	1.489	1.340
99) m,p-xylene	0.570	0.488	0.586	0.609	0.612	0.656	0.614	0.627	0.585	0.396
100) o-xylene	0.584	0.505	0.600	0.639	0.644	0.649	0.661	0.655	0.587	0.342
101) styrene	0.927	0.763	0.936	1.044	1.047	1.063	1.052	1.063	0.989	0.987
102) Butyl Acrylate									0.000#	-1.00
103) bromoform	0.275	0.239	0.271	0.309	0.327	0.280	0.339	0.324	0.294	0.295
104) I 1,4-dichlorobenzene-d										ISTD-----
105) isopropylbenzene	2.411	2.181	2.464	2.679	2.795	2.847	2.726	2.678	2.474	1.835
106) 4-bromofluorobenzene (s)	0.783	0.766	0.765	0.776	0.784	0.787	0.734	0.762	0.770	0.766
107) bromobenzene	0.670	0.686	0.732	0.759	0.777	0.810	0.732	0.769	0.729	0.741
108) cyclohexanone ** This compound does not meet Initial Calibration criteria	0.080	0.034	0.096		0.093	0.102	0.093		0.083	30.12
										----- Linear regression ----- Coefficient = 0.9708
										Response Ratio = -0.10526 + 0.09442 *A
109) 1,1,2,2-tetrachloroethane	0.701	0.695	0.693	0.718	0.770	0.813	0.755	0.764	0.720	0.633
110) trans-1,4-dichloro-2-butene	0.167	0.137	0.158	0.173	0.153	0.171	0.171	0.155	0.161	7.72
111) 1,2,3-trichloropropane	0.162	0.141	0.164	0.176	0.172	0.170	0.174	0.168	0.166	6.78

Initial Calibration Summary

Page 5 of 5

Job Number: JC1106

Sample: V3D4810-ICC4810
Lab FileID: 3D112242.DAccount: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

112)	n-propylbenzene	2.696 2.328 2.834 2.877 2.973 3.167 2.773 2.928 2.756 1.986	2.732	12.45
113)	4-Ethyltoluene		0.000#	-1.00
114)	2-chlorotoluene	0.610 0.588 0.616 0.654 0.688 0.649 0.648 0.669 0.621	0.638	4.96
115)	4-chlorotoluene	0.612 0.559 0.576 0.662 0.669 0.712 0.637 0.664 0.636 0.468	0.620	11.27
116)	1,3,5-trimethylbenzene	2.069 1.885 2.090 2.272 2.399 2.383 2.357 2.276 2.110 1.510	2.135	12.87
117)	tert-butylbenzene	1.609 1.377 1.752 1.916 2.058 1.903 2.151 1.882 1.690 1.254	1.759	16.21
118)	pentachloroethane	0.429 0.328 0.438 0.508 0.576 0.510 0.617 0.483 0.430	0.480	17.96
119)	1,2,4-trimethylbenzene	2.105 1.826 2.055 2.249 2.360 2.428 2.278 2.277 2.131 1.728	2.144	10.52
120)	sec-butylbenzene	2.775 2.195 2.865 3.045 3.242 3.249 3.223 3.084 2.812 1.625	2.811	18.62
121)	1,3-dichlorobenzene	1.321 1.235 1.347 1.462 1.486 1.526 1.472 1.508 1.405 1.008	1.377	11.56
122)	p-isopropyltoluene	2.376 2.151 2.421 2.598 2.760 2.779 2.700 2.650 2.431 1.683	2.455	13.66
123)	1,4-dichlorobenzene	1.402 1.334 1.415 1.498 1.515 1.749 1.521 1.538 1.477 1.324	1.477	8.29
124)	1,2-dichlorobenzene	1.441 1.263 1.489 1.557 1.589 1.672 1.610 1.585 1.491 1.149	1.484	11.02
125)	1,4-Diethylbenzene		0.000#	-1.00
126)	n-butylbenzene	1.206 0.992 1.278 1.333 1.367 1.386 1.339 1.373 1.287	1.285	9.63
127)	1,2,4,5-Tetramethylbenzene		0.000#	-1.00
128)	1,2-dibromo-3-chloropropane	0.195 0.192 0.214 0.240 0.208 0.259 0.228 0.201	0.217	10.80
129)	1,3,5-trichlorobenzene	1.477 1.359 1.600 1.694 1.748 1.711 1.869 1.739 1.581 0.984	1.576	16.14
130)	1,2,4-trichlorobenzene	1.499 1.264 1.531 1.747 1.828 1.712 1.936 1.753 1.633 1.299	1.620	13.61
131)	hexachlorobutadiene	0.761 0.556 0.826 0.793 0.859 0.898 0.933 0.811 0.781	0.802	13.46
132)	naphthalene	3.116 2.773 3.113 3.521 3.857 3.459 4.038 3.662 3.368 3.411	3.432	10.86
133)	1,2,3-trichlorobenzene	1.567 1.458 1.664 1.776 1.877 1.825 1.916 1.795 1.647 1.295	1.682	11.72
134)	hexachloroethane	0.417 0.414 0.537 0.616 0.460 0.675 0.500 0.437	0.507	19.00
135)	Benzyl chloride	1.391 1.506 1.581 1.359 1.406 1.711 1.450 1.363 1.315 1.343	1.442	8.61

(##) = Out of Range ### Number of calibration levels exceeded format ###

M3D4810.M

Thu Aug 06 13:51:48 2015

3D

Initial Calibration Verification

Page 1 of 3

Job Number: JC1106
Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample: V3D4810-ICV4810
Lab FileID: 3D112247.D

Evaluate Continuing Calibration Report

Data File : C:\msdchem\1\DATA\v3d4801-4810\3d112247.d Vial: 32
 Acq On : 30 Jul 2015 1:41 am Operator: ximenac
 Sample : ICV4810-50 Inst : MS3D
 Misc : MS88759, V3D4810, 5, , , , 1 Multiplr: 1.00
 MS Integration Params: rteint.p

Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)
 Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 Last Update : Thu Aug 06 11:10:54 2015
 Response via : Multiple Level Calibration

Min. RRF : 0.010 Min. Rel. Area : 50% Max. R.T. Dev 0.30min
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)	R.T.
1 I	Tert Butyl Alcohol-d9	1.000	1.000	0.0	102	0.00	7.18
2	1,4-dioxane	0.081	0.089	-9.9	103	0.00	11.04
3 M	tertiary butyl alcohol	1.080	1.088	-0.7	103	-0.01	7.29
4 I	pentafluorobenzene	1.000	1.000	0.0	97	0.00	9.39
5	FREON 115			-----NA-----			
6	FREON 152A			-----NA-----			
7	chlorotrifluoroethene			-----NA-----			
8	chlorodifluoromethane	0.679	0.568	16.3	81	0.00	3.85
9	dichlorodifluoromethane	0.909	0.674	25.9	72	0.00	3.85
10	chloromethane	0.723	0.600	17.0	86	0.00	4.16
11	vinyl chloride	0.725	0.661	8.8	85	0.00	4.41
12	bromomethane	0.468	0.433	7.5	93	0.00	5.05
13	chloroethane	0.307	0.316	-2.9	104	0.00	5.21
14	1,3-butadiene			-----NA-----			
15	vinyl bromide			-----NA-----			
16	trichlorofluoromethane	0.756	0.656	13.2	81	0.00	5.68
17	pentane			-----NA-----			
18	ethyl ether	0.187	0.176	5.9	96	0.00	6.07
19	freon 123a			-----NA-----			
20	FREON 123			-----NA-----			
21	acrolein	0.075	0.076	-1.3	108	0.00	6.32
22	1,1-dichloroethene	0.595	0.499	16.1	78	0.00	6.49
23	acetone	0.112	0.108	3.6	110	0.00	6.54
24	allyl chloride	0.185	0.199	-7.6	96	0.00	7.02
25	acetonitrile	0.037	0.038	-2.7	104	0.00	6.97
26 m	acetaldehyde			-----NA-----			
27	iodomethane	0.846	0.822	2.8	87	0.00	6.76
28	iso-butyl alcohol	0.012	0.012	0.0	101	0.00	9.68
29	carbon disulfide	1.504	1.353	10.0	84	0.00	6.89
30	methylene chloride	0.446	0.429	3.8	91	0.00	7.20
31	methyl acetate	0.249	0.258	-3.6	99	0.00	7.00
32	methyl tert butyl ether	1.320	1.295	1.9	95	0.00	7.53
33	trans-1,2-dichloroethene	0.543	0.492	9.4	85	0.00	7.58
34	di-isopropyl ether	1.396	1.371	1.8	97	0.00	8.12
35	ethyl tert-butyl ether	1.315	1.405	-6.8	99	0.00	8.59
36	2-butanone	0.036	0.037	-2.8	105	0.00	8.84
37 M	1,1-dichloroethane	0.645	0.638	1.1	91	0.00	8.14
38	chloroprene	0.472	0.469	0.6	92	0.00	8.25
39	acrylonitrile	0.129	0.137	-6.2	99	0.00	7.52
40	vinyl acetate	0.048	0.055	-14.6	107	0.00	8.13
41	ethyl acetate	0.049	0.047	4.1	107	0.00	8.86

6.7.2

6

Initial Calibration Verification

Page 2 of 3

Job Number: JC1106

Sample: V3D4810-ICV4810
Lab FileID: 3D112247.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

42	2,2-dichloropropane	0.675	0.557	17.5	84	0.00	8.88
43	cis-1,2-dichloroethene	0.436	0.391	10.3	91	0.00	8.88
44	propionitrile	0.048	0.047	2.1	97	0.00	8.94
45	methyl acrylate	0.324	0.335	-3.4	99	0.00	8.95
46	bromochloromethane	0.210	0.213	-1.4	95	0.00	9.19
47	tetrahydrofuran	0.126	0.120	4.8	103	0.00	9.22
48	chloroform	0.622	0.604	2.9	92	0.00	9.24
49	Tert-Butyl Formate	0.362	0.396	-9.4	103	0.00	9.28
50 S	dibromofluoromethane (s)	0.383	0.388	-1.3	97	0.00	9.44
51 S	1,2-dichloroethane-d4 (s)	0.355	0.346	2.5	96	0.00	9.86
52	freon 113	0.366	0.344	6.0	89	0.00	6.45
53	methacrylonitrile	0.188	0.191	-1.6	102	0.00	9.12
54	1,1,1-trichloroethane	0.642	0.594	7.5	88	0.00	9.50
55	tert amyl alcohol			-----NA-----			
56	2,2,4-Trimethylpentane			-----NA-----			
57	tert-amyl methyl ether	1.314	1.315	-0.1	100	0.00	9.97
58 I	1,4-difluorobenzene	1.000	1.000	0.0	96	0.00	10.31
59	epichlorohydrin	0.027	0.028	-3.7	109	0.00	11.57
60	n-butyl alcohol	0.008	0.009#	-12.5	104	0.00	10.44
61 M	cyclohexane	0.586	0.512	12.6	83	0.00	9.57
62	carbon tetrachloride	0.484	0.472	2.5	85	0.00	9.71
63	1,1-dichloropropene	0.360	0.351	2.5	87	0.00	9.68
64	hexane	0.322	0.210	34.8#	62	0.00	7.87
65 M	benzene	1.175	1.123	4.4	90	0.00	9.94
66	heptane	0.183	0.163	10.9	86	0.00	10.10
67	isopropyl acetate	0.511	0.624	-22.1	116	0.00	9.86
68	1,2-dichloroethane	0.351	0.362	-3.1	96	0.00	9.95
69	trichloroethene	0.284	0.267	6.0	89	0.00	10.66
70	ethyl acrylate			-----NA-----			
71	tert amyl ethyl ether			-----NA-----			
72	2-nitropropane	0.104	0.108	-3.8	106	0.00	11.43
73	2-chloroethyl vinyl ether	0.180	0.217	-20.6	114	0.00	11.45
74	methyl methacrylate	0.402	0.440	-9.5	102	0.00	10.92
75	1,2-dichloropropane	0.307	0.308	-0.3	94	0.00	10.92
76	methylcyclohexane	0.531	0.545	-2.6	88	0.00	10.87
77	dibromomethane	0.181	0.185	-2.2	95	0.00	11.08
78	bromodichloromethane	0.374	0.379	-1.3	94	0.00	11.21
79	cis-1,3-dichloropropene	0.439	0.462	-5.2	95	0.00	11.67
80 S	toluene-d8 (s)	1.095	1.102	-0.6	96	0.00	11.96
81	4-methyl-2-pentanone	0.112	0.120	-7.1	104	0.00	11.76
82	toluene	0.646	0.638	1.2	88	0.00	12.03
83	3-methyl-1-butanol	0.008	0.008#	0.0	104	0.00	11.79
84	trans-1,3-dichloropropene	0.390	0.397	-1.8	91	0.00	12.23
85	ethyl methacrylate	0.345	0.348	-0.9	95	0.00	12.23
86	1,1,2-trichloroethane	0.217	0.218	-0.5	96	0.00	12.45
87	2-hexanone	0.091	0.096	-5.5	104	0.00	12.62
88 I	chlorobenzene-d5	1.000	1.000	0.0	97	0.00	13.48
89	tetrachloroethene	0.426	0.445	-4.5	98	0.00	12.63
90	1,3-dichloropropane	0.485	0.483	0.4	95	0.00	12.63
91	butyl acetate	0.201	0.221	-10.0	105	0.00	12.69
92	3,3-Dimethyl-1-Butanol	0.048	0.041	14.6	99	0.00	12.79
93	dibromochloromethane	0.378	0.398	-5.3	96	0.00	12.89
94	1,2-dibromoethane	0.319	0.331	-3.8	97	0.00	13.05
95	n-Butyl Ether	1.513	1.413	6.6	91	0.00	13.42
96	chlorobenzene	0.897	0.916	-2.1	95	0.00	13.51
97	1,1,1,2-tetrachloroethane	0.398	0.403	-1.3	91	0.00	13.56
98	ethylbenzene	1.536	1.457	5.1	90	0.00	13.56
99	m,p-xylene	0.574	0.572	0.3	91	0.00	13.67

Initial Calibration Verification

Page 3 of 3

Job Number: JC1106

Sample: V3D4810-ICV4810
Lab FileID: 3D112247.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

100	o-xylene	0.587	0.610	-3.9	93	0.00	14.08
101	styrene	0.987	1.023	-3.6	95	0.00	14.09
102	Butyl Acrylate			NA			
103	bromoform	0.295	0.323	-9.5	101	0.00	14.36
104 I	1,4-dichlorobenzene-d4	1.000	1.000	0.0	101	0.00	15.79
105	isopropylbenzene	2.509	2.337	6.9	88	0.00	14.42
106 S	4-bromofluorobenzene (s)	0.769	0.763	0.8	100	0.00	14.63
107	bromobenzene	0.741	0.726	2.0	97	0.00	14.82
-----		True	Calc.	% Drift	-----		
108	cyclohexanone	500.000	487.845	2.4	243	0.00	14.59
-----		AvgRF	CCRF	% Dev	-----		
109	1,1,2,2-tetrachloroethane	0.726	0.705	2.9	100	0.00	14.73
110	trans-1,4-dichloro-2-bute	0.161	0.158	1.9	101	0.00	14.77
111	1,2,3-trichloropropane	0.166	0.164	1.2	101	0.00	14.80
112	n-propylbenzene	2.732	2.714	0.7	96	0.00	14.83
113	4-Ethyltoluene		NA				
114	2-chlorotoluene	0.638	0.607	4.9	94	0.00	14.98
115	4-chlorotoluene	0.620	0.623	-0.5	95	0.00	15.08
116	1,3,5-trimethylbenzene	2.135	2.034	4.7	91	0.00	14.99
117	tert-butylbenzene	1.759	1.683	4.3	89	0.00	15.33
118	pentachloroethane	0.480	0.402	16.2	80	0.00	15.42
119	1,2,4-trimethylbenzene	2.144	2.167	-1.1	98	0.00	15.38
120	sec-butylbenzene	2.811	2.657	5.5	89	0.00	15.55
121	1,3-dichlorobenzene	1.377	1.409	-2.3	98	0.00	15.73
122	p-isopropyltoluene	2.455	2.367	3.6	92	0.00	15.66
123	1,4-dichlorobenzene	1.477	1.465	0.8	99	0.00	15.81
124	1,2-dichlorobenzene	1.484	1.554	-4.7	101	0.00	16.19
125	1,4-Diethylbenzene		NA				
126	n-butylbenzene	1.285	1.237	3.7	94	0.00	16.07
127	1,2,4,5-Tetramethylbenzen		NA				
128	1,2-dibromo-3-chloropropa	0.217	0.226	-4.1	107	0.00	16.93
129	1,3,5-trichlorobenzene	1.576	1.641	-4.1	98	0.00	17.09
130	1,2,4-trichlorobenzene	1.620	1.699	-4.9	99	0.00	17.66
131	hexachlorobutadiene	0.802	0.717	10.6	92	0.00	17.76
132	naphthalene	3.432	3.555	-3.6	102	0.00	17.90
133	1,2,3-trichlorobenzene	1.682	1.724	-2.5	98	0.00	18.11
134	hexachloroethane	0.507	0.479	5.5	90	0.00	16.44
135	Benzyl chloride	1.442	1.418	1.7	106	0.00	15.92

(#) = Out of Range
3D112242.D M3D4810.M

SPCC's out = 0 CCC's out = 0
Thu Aug 06 13:51:40 2015 3D

6.7.2
6

Continuing Calibration Summary

Page 1 of 3

Job Number: JC1106

Sample: V3D4822-CC4810

Account: UTC United Technologies Corporation

Lab FileID: 3D112517.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Evaluate Continuing Calibration Report

Data File : C:\msdchem\1\DATA\3D112517.D

Vial: 5

Acq On : 11 Aug 2015 11:16 am

Operator: ximenac

Sample : cc4810-20

Inst : MS3D

Misc : MS89341,V3D4822,5,,,1

Multiplr: 1.00

MS Integration Params: rteint.p

Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)

Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

Last Update : Thu Aug 06 11:10:54 2015

Response via : Multiple Level Calibration

Min. RRF : 0.010 Min. Rel. Area : 50% Max. R.T. Dev 0.30min

Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)	R.T.
1 I	Tert Butyl Alcohol-d9	1.000	1.000	0.0	110	-0.02	7.16
2	1,4-dioxane	0.081	0.093	-14.8	116	0.00	11.04
3 M	tertiary butyl alcohol	1.080	1.074	0.6	106	0.00	7.30
4 I	pentafluorobenzene	1.000	1.000	0.0	100	0.00	9.39
5	FREON 115			-----NA-----			
6	FREON 152A			-----NA-----			
7	chlorotrifluoroethene			-----NA-----			
8	chlorodifluoromethane	0.679	0.664	2.2	96	0.00	3.86
9	dichlorodifluoromethane	0.909	0.742	18.4	78	0.01	3.86
10	chloromethane	0.723	0.551	23.8#	85	0.00	4.16
11	vinyl chloride	0.725	0.630	13.1	88	0.00	4.41
12	bromomethane	0.468	0.407	13.0	90	0.00	5.05
13	chloroethane	0.307	0.276	10.1	95	0.01	5.22
14	1,3-butadiene			-----NA-----			
15	vinyl bromide			-----NA-----			
16	trichlorofluoromethane	0.756	0.773	-2.2	99	0.01	5.68
17	pentane			-----NA-----			
18	ethyl ether	0.187	0.197	-5.3	104	0.00	6.07
19	freon 123a			-----NA-----			
20	FREON 123			-----NA-----			
21	acrolein	0.075	0.071	5.3	96	0.00	6.32
22	1,1-dichloroethene	0.595	0.602	-1.2	98	0.00	6.49
23	acetone	0.112	0.137	-22.3#	113	0.00	6.54
24	allyl chloride	0.185	0.197	-6.5	97	0.00	7.02
25	acetonitrile	0.037	0.043	-16.2	115	0.00	6.98
26 m	acetaldehyde			-----NA-----			
27	iodomethane	0.846	0.944	-11.6	104	0.00	6.76
28	iso-butyl alcohol	0.012	0.013	-8.3	105	0.00	9.69
29	carbon disulfide	1.504	1.599	-6.3	101	0.00	6.90
30	methylene chloride	0.446	0.492	-10.3	107	0.00	7.20
31	methyl acetate	0.249	0.276	-10.8	109	0.00	6.99
32	methyl tert butyl ether	1.320	1.387	-5.1	101	0.00	7.53
33	trans-1,2-dichloroethene	0.543	0.577	-6.3	101	0.00	7.58
34	di-isopropyl ether	1.396	1.250	10.5	93	0.00	8.12
35	ethyl tert-butyl ether	1.315	1.276	3.0	94	0.00	8.59
36	2-butanone	0.036	0.037	-2.8	92	0.00	8.85
37 M	1,1-dichloroethane	0.645	0.700	-8.5	100	0.00	8.14
38	chloroprene	0.472	0.404	14.4	86	0.00	8.25
39	acrylonitrile	0.129	0.151	-17.1	105	0.00	7.53
40	vinyl acetate	0.048	0.050	-4.2	99	0.00	8.13
41	ethyl acetate	0.049	0.044	10.2	85	0.00	8.87

Continuing Calibration Summary

Page 2 of 3

Job Number: JC1106

Sample: V3D4822-CC4810
Lab FileID: 3D112517.D

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

42	2,2-dichloropropane	0.675	0.716	-6.1	108	0.00	8.89
43	cis-1,2-dichloroethene	0.436	0.421	3.4	97	0.00	8.88
44	propionitrile	0.048	0.055	-14.6	108	0.00	8.94
45	methyl acrylate	0.324	0.358	-10.5	98	0.00	8.95
46	bromochloromethane	0.210	0.231	-10.0	101	0.00	9.19
47	tetrahydrofuran	0.126	0.122	3.2	93	0.00	9.22
48	chloroform	0.622	0.648	-4.2	101	0.00	9.24
49	Tert-Butyl Formate	0.362	0.346	4.4	95	0.00	9.28
50 S	dibromofluoromethane (s)	0.383	0.412	-7.6	108	0.00	9.44
51 S	1,2-dichloroethane-d4 (s)	0.355	0.381	-7.3	107	0.00	9.86
52	freon 113	0.366	0.371	-1.4	101	0.01	6.46
53	methacrylonitrile	0.188	0.184	2.1	95	0.00	9.13
54	1,1,1-trichloroethane	0.642	0.714	-11.2	107	0.00	9.50
55	tert amyl alcohol			-----NA-----			
56	2,2,4-Trimethylpentane			-----NA-----			
57	tert-amyl methyl ether	1.314	1.243	5.4	98	0.00	9.97
58 I	1,4-difluorobenzene	1.000	1.000	0.0	101	0.00	10.31
59	epichlorohydrin	0.027	0.026	3.7	99	0.00	11.58
60	n-butyl alcohol	0.008	0.008#	0.0	103	0.00	10.44
61 M	cyclohexane	0.586	0.582	0.7	101	0.00	9.57
62	carbon tetrachloride	0.484	0.562	-16.1	109	0.00	9.71
63	1,1-dichloropropene	0.360	0.370	-2.8	96	0.00	9.67
64	hexane	0.322	0.271	15.8	89	0.00	7.87
65 M	benzene	1.175	1.196	-1.8	100	0.00	9.94
66	heptane	0.183	0.172	6.0	97	0.00	10.11
67	isopropyl acetate	0.511	0.494	3.3	95	0.00	9.86
68	1,2-dichloroethane	0.351	0.385	-9.7	104	0.00	9.96
69	trichloroethene	0.284	0.286	-0.7	100	0.00	10.66
70	ethyl acrylate			-----NA-----			
71	tert amyl ethyl ether			-----NA-----			
72	2-nitropropane	0.104	0.100	3.8	88	0.01	11.44
73	2-chloroethyl vinyl ether	0.180	0.181	-0.6	96	0.00	11.45
74	methyl methacrylate	0.402	0.408	-1.5	93	0.00	10.93
75	1,2-dichloropropane	0.307	0.316	-2.9	101	0.00	10.92
76	methylcyclohexane	0.531	0.521	1.9	95	0.00	10.88
77	dibromomethane	0.181	0.207	-14.4	105	0.00	11.08
78	bromodichloromethane	0.374	0.398	-6.4	102	0.00	11.21
79	cis-1,3-dichloropropene	0.439	0.461	-5.0	95	0.00	11.67
80 S	toluene-d8 (s)	1.095	1.102	-0.6	101	0.00	11.96
81	4-methyl-2-pentanone	0.112	0.119	-6.2	100	0.00	11.77
82	toluene	0.646	0.678	-5.0	95	0.00	12.04
83	3-methyl-1-butanol	0.008	0.008#	0.0	102	0.00	11.79
84	trans-1,3-dichloropropene	0.390	0.417	-6.9	97	0.00	12.24
85	ethyl methacrylate	0.345	0.314	9.0	86	0.00	12.23
86	1,1,2-trichloroethane	0.217	0.232	-6.9	102	0.00	12.45
87	2-hexanone	0.091	0.085	6.6	86	0.00	12.63
88 I	chlorobenzene-d5	1.000	1.000	0.0	102	0.00	13.48
89	tetrachloroethene	0.426	0.429	-0.7	95	0.00	12.63
90	1,3-dichloropropane	0.485	0.512	-5.6	101	0.00	12.63
91	butyl acetate	0.201	0.187	7.0	95	0.00	12.70
92	3,3-Dimethyl-1-Butanol	0.048	0.037	22.9#	91	0.00	12.79
93	dibromochloromethane	0.378	0.422	-11.6	104	0.00	12.90
94	1,2-dibromoethane	0.319	0.345	-8.2	99	0.00	13.05
95	n-Butyl Ether	1.513	1.278	15.5	84	0.00	13.42
96	chlorobenzene	0.897	0.948	-5.7	99	0.00	13.51
97	1,1,2-tetrachloroethane	0.398	0.434	-9.0	105	0.00	13.56
98	ethylbenzene	1.536	1.509	1.8	96	0.00	13.56
99	m,p-xylene	0.574	0.585	-1.9	96	0.00	13.67

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Continuing Calibration Summary

Page 3 of 3

Job Number: JC1106

Sample: V3D4822-CC4810
Lab FileID: 3D112517.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

100	o-xylene	0.587	0.606	-3.2	95	0.00	14.09
101	styrene	0.987	0.933	5.5	90	0.00	14.10
102	Butyl Acrylate			-----NA-----			
103	bromoform	0.295	0.329	-11.5	104	0.00	14.36
104 I	1,4-dichlorobenzene-d4	1.000	1.000	0.0	100	0.00	15.79
105	isopropylbenzene	2.509	2.476	1.3	93	0.00	14.43
106 S	4-bromofluorobenzene (s)	0.769	0.773	-0.5	102	0.00	14.63
107	bromobenzene	0.741	0.755	-1.9	99	0.00	14.83
		-----True	Calc.	% Drift			
108	cyclohexanone	200.000	259.909	-30.0#	95	0.00	14.59
		-----AvgRF	CCRF	% Dev			
109	1,1,2,2-tetrachloroethane	0.726	0.805	-10.9	106	0.00	14.73
110	trans-1,4-dichloro-2-bute	0.161	0.148	8.1	87	0.00	14.78
111	1,2,3-trichloropropane	0.166	0.195	-17.5	112	0.00	14.81
112	n-propylbenzene	2.732	2.792	-2.2	96	0.00	14.84
113	4-Ethyltoluene			-----NA-----			
114	2-chlorotoluene	0.638	0.655	-2.7	98	0.00	14.99
115	4-chlorotoluene	0.620	0.650	-4.8	98	0.00	15.08
116	1,3,5-trimethylbenzene	2.135	2.215	-3.7	98	0.00	14.99
117	tert-butylbenzene	1.759	1.695	3.6	90	0.00	15.33
118	pentachloroethane	0.480	0.518	-7.9	108	0.00	15.42
119	1,2,4-trimethylbenzene	2.144	2.170	-1.2	96	0.00	15.38
120	sec-butylbenzene	2.811	2.862	-1.8	93	0.00	15.55
121	1,3-dichlorobenzene	1.377	1.470	-6.8	98	0.00	15.74
122	p-isopropyltoluene	2.455	2.469	-0.6	94	0.00	15.67
123	1,4-dichlorobenzene	1.477	1.537	-4.1	100	0.00	15.82
124	1,2-dichlorobenzene	1.484	1.600	-7.8	101	0.00	16.20
125	1,4-Diethylbenzene			-----NA-----			
126	n-butylbenzene	1.285	1.296	-0.9	95	0.00	16.07
127	1,2,4,5-Tetramethylbenzen			-----NA-----			
128	1,2-dibromo-3-chloropropa	0.217	0.233	-7.4	103	0.00	16.93
129	1,3,5-trichlorobenzene	1.576	1.709	-8.4	99	0.00	17.10
130	1,2,4-trichlorobenzene	1.620	1.789	-10.4	102	0.00	17.66
131	hexachlorobutadiene	0.802	0.806	-0.5	100	0.00	17.76
132	naphthalene	3.432	3.789	-10.4	104	0.00	17.90
133	1,2,3-trichlorobenzene	1.682	1.893	-12.5	106	0.00	18.12
134	hexachloroethane	0.507	0.503	0.8	101	0.00	16.44
135	Benzyl chloride	1.442	1.443	-0.1	106	0.00	15.93

(#) = Out of Range
3D112241.D M3D4810.M

SPCC's out = 0 CCC's out = 0
Tue Aug 11 16:10:17 2015 3D

Continuing Calibration Summary

Page 1 of 3

Job Number: JC1106
Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample: V3D4823-CC4810
Lab FileID: 3D112542.D

Evaluate Continuing Calibration Report

Data File : C:\msdchem\1\DATA\3D\v3d4822\3D112542.D Vial: 35
 Acq On : 11 Aug 2015 11:17 pm Operator: ximenac
 Sample : cc4810-50 Inst : MS3D
 Misc : MS89468, V3D4823, 5, , , , 1 Multiplr: 1.00
 MS Integration Params: rteint.p

Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)
 Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 Last Update : Thu Aug 06 11:10:54 2015
 Response via : Multiple Level Calibration

Min. RRF : 0.010 Min. Rel. Area : 50% Max. R.T. Dev 0.30min
 Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)	R.T.
1 I	Tert Butyl Alcohol-d9	1.000	1.000	0.0	94	0.00	7.17
2	1,4-dioxane	0.081	0.096	-18.5	101	0.00	11.04
3 M	tertiary butyl alcohol	1.080	1.101	-1.9	96	-0.02	7.28
4 I	pentafluorobenzene	1.000	1.000	0.0	89	0.00	9.39
5	FREON 115			-----NA-----			
6	FREON 152A			-----NA-----			
7	chlorotrifluoroethene			-----NA-----			
8	chlorodifluoromethane	0.679	0.778	-14.6	102	0.00	3.85
9	dichlorodifluoromethane	0.909	1.047	-15.2	103	-0.01	3.84
10	chloromethane	0.723	0.802	-10.9	105	0.00	4.16
11	vinyl chloride	0.725	0.922	-27.2#	109	0.00	4.42
12	bromomethane	0.468	0.548	-17.1	109	0.00	5.04
13	chloroethane	0.307	0.358	-16.6	109	0.00	5.21
14	1,3-butadiene			-----NA-----			
15	vinyl bromide			-----NA-----			
16	trichlorofluoromethane	0.756	0.995	-31.6#	113	0.00	5.67
17	pentane			-----NA-----			
18	ethyl ether	0.187	0.197	-5.3	99	0.00	6.07
19	freon 123a			-----NA-----			
20	FREON 123			-----NA-----			
21	acrolein	0.075	0.075	0.0	98	0.00	6.32
22	1,1-dichloroethene	0.595	0.687	-15.5	99	0.00	6.48
23	acetone	0.112	0.122	-8.9	114	0.00	6.54
24	allyl chloride	0.185	0.206	-11.4	91	0.00	7.01
25	acetonitrile	0.037	0.042	-13.5	106	-0.01	6.97
26 m	acetaldehyde			-----NA-----			
27	iodomethane	0.846	1.016	-20.1#	99	0.00	6.76
28	iso-butyl alcohol	0.012	0.013	-8.3	98	0.00	9.68
29	carbon disulfide	1.504	1.800	-19.7	103	0.00	6.89
30	methylene chloride	0.446	0.513	-15.0	101	0.00	7.20
31	methyl acetate	0.249	0.278	-11.6	98	0.00	6.99
32	methyl tert butyl ether	1.320	1.440	-9.1	97	0.00	7.53
33	trans-1,2-dichloroethene	0.543	0.624	-14.9	99	0.00	7.57
34	di-isopropyl ether	1.396	1.415	-1.4	92	0.00	8.12
35	ethyl tert-butyl ether	1.315	1.423	-8.2	92	0.00	8.58
36	2-butanone	0.036	0.038	-5.6	98	0.00	8.85
37 M	1,1-dichloroethane	0.645	0.752	-16.6	98	0.00	8.14
38	chloroprene	0.472	0.494	-4.7	89	0.00	8.25
39	acrylonitrile	0.129	0.150	-16.3	100	0.00	7.52
40	vinyl acetate	0.048	0.054	-12.5	95	0.00	8.13
41	ethyl acetate	0.049	0.045	8.2	93	0.00	8.86

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Continuing Calibration Summary

Page 2 of 3

Job Number: JC1106

Sample: V3D4823-CC4810
Lab FileID: 3D112542.D

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

42	2,2-dichloropropane	0.675	0.728	-7.9	101	0.00	8.88
43	cis-1,2-dichloroethene	0.436	0.457	-4.8	98	0.00	8.88
44	propionitrile	0.048	0.054	-12.5	102	0.00	8.93
45	methyl acrylate	0.324	0.358	-10.5	97	0.00	8.94
46	bromochloromethane	0.210	0.246	-17.1	101	0.00	9.19
47	tetrahydrofuran	0.126	0.124	1.6	97	0.00	9.23
48	chloroform	0.622	0.712	-14.5	99	0.00	9.24
49	Tert-Butyl Formate	0.362	0.379	-4.7	91	0.00	9.27
50 S	dibromofluoromethane (s)	0.383	0.417	-8.9	96	0.00	9.44
51 S	1,2-dichloroethane-d4 (s)	0.355	0.388	-9.3	99	0.00	9.86
52	freon 113	0.366	0.453	-23.8#	108	0.00	6.45
53	methacrylonitrile	0.188	0.187	0.5	92	0.00	9.12
54	1,1,1-trichloroethane	0.642	0.787	-22.6#	107	0.00	9.50
55	tert amyl alcohol			-----NA-----			
56	2,2,4-Trimethylpentane			-----NA-----			
57	tert-amyl methyl ether	1.314	1.401	-6.6	98	0.00	9.97
58 I	1,4-difluorobenzene	1.000	1.000	0.0	91	0.00	10.31
59	epichlorohydrin	0.027	0.027	0.0	103	0.00	11.57
60	n-butyl alcohol	0.008	0.009#	-12.5	98	0.00	10.44
61 M	cyclohexane	0.586	0.661	-12.8	102	0.00	9.57
62	carbon tetrachloride	0.484	0.636	-31.4#	109	0.00	9.71
63	1,1-dichloropropene	0.360	0.423	-17.5	100	0.00	9.67
64	hexane	0.322	0.317	1.6	90	0.00	7.87
65 M	benzene	1.175	1.325	-12.8	101	0.00	9.94
66	heptane	0.183	0.200	-9.3	101	0.00	10.11
67	isopropyl acetate	0.511	0.537	-5.1	95	0.00	9.86
68	1,2-dichloroethane	0.351	0.408	-16.2	102	0.00	9.95
69	trichloroethene	0.284	0.322	-13.4	102	0.00	10.66
70	ethyl acrylate			-----NA-----			
71	tert amyl ethyl ether			-----NA-----			
72	2-nitropropane	0.104	0.106	-1.9	100	0.00	11.44
73	2-chloroethyl vinyl ether	0.180	0.197	-9.4	98	0.00	11.45
74	methyl methacrylate	0.402	0.429	-6.7	94	0.00	10.92
75	1,2-dichloropropane	0.307	0.345	-12.4	101	0.00	10.92
76	methylcyclohexane	0.531	0.654	-23.2#	100	0.00	10.87
77	dibromomethane	0.181	0.222	-22.7#	108	0.00	11.08
78	bromodichloromethane	0.374	0.436	-16.6	103	0.00	11.21
79	cis-1,3-dichloropropene	0.439	0.498	-13.4	97	0.00	11.67
80 S	toluene-d8 (s)	1.095	1.134	-3.6	94	0.00	11.96
81	4-methyl-2-pentanone	0.112	0.119	-6.2	98	0.00	11.76
82	toluene	0.646	0.759	-17.5	100	0.00	12.03
83	3-methyl-1-butanol	0.008	0.009#	-12.5	104	0.00	11.78
84	trans-1,3-dichloropropene	0.390	0.457	-17.2	100	0.00	12.23
85	ethyl methacrylate	0.345	0.339	1.7	88	0.00	12.23
86	1,1,2-trichloroethane	0.217	0.247	-13.8	103	0.00	12.45
87	2-hexanone	0.091	0.085	6.6	88	0.00	12.62
88 I	chlorobenzene-d5	1.000	1.000	0.0	95	0.00	13.48
89	tetrachloroethene	0.426	0.453	-6.3	97	0.00	12.62
90	1,3-dichloropropane	0.485	0.531	-9.5	101	0.00	12.63
91	butyl acetate	0.201	0.206	-2.5	96	0.00	12.69
92	3,3-Dimethyl-1-Butanol	0.048	0.041	14.6	95	0.00	12.79
93	dibromochloromethane	0.378	0.448	-18.5	105	0.00	12.89
94	1,2-dibromoethane	0.319	0.355	-11.3	101	0.00	13.05
95	n-Butyl Ether	1.513	1.441	4.8	90	0.00	13.42
96	chlorobenzene	0.897	1.016	-13.3	102	0.00	13.51
97	1,1,1,2-tetrachloroethane	0.398	0.476	-19.6	105	0.00	13.56
98	ethylbenzene	1.536	1.663	-8.3	101	0.00	13.56
99	m,p-xylene	0.574	0.647	-12.7	101	0.00	13.66

Continuing Calibration Summary

Page 3 of 3

Job Number: JC1106

Sample: V3D4823-CC4810

Account: UTC United Technologies Corporation

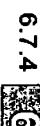
Lab FileID: 3D112542.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

100	o-xylene	0.587	0.667	-13.6	99	0.00	14.08
101	styrene	0.987	1.041	-5.5	94	0.00	14.09
102	Butyl Acrylate			-----NA-----			
103	bromoform	0.295	0.337	-14.2	103	0.00	14.36
104 I	1,4-dichlorobenzene-d4	1.000	1.000	0.0	93	0.00	15.79
105	isopropylbenzene	2.509	2.882	-14.9	100	0.00	14.42
106 S	4-bromofluorobenzene (s)	0.769	0.776	-0.9	93	0.00	14.63
107	bromobenzene	0.741	0.817	-10.3	100	0.00	14.82
-----		True	Calc.	% Drift	-----		
108	cyclohexanone	500.000	257.854	48.4#	104	0.00	14.59
-----		AvgRF	CCRF	% Dev	-----		
109	1,1,2,2-tetrachloroethane	0.726	0.814	-12.1	105	0.00	14.73
110	trans-1,4-dichloro-2-bute	0.161	0.134	16.8	79	0.00	14.77
111	1,2,3-trichloropropane	0.166	0.193	-16.3	110	0.00	14.80
112	n-propylbenzene	2.732	3.123	-14.3	101	0.00	14.83
113	4-Ethyltoluene			-----NA-----			
114	2-chlorotoluene	0.638	0.729	-14.3	104	0.00	14.98
115	4-chlorotoluene	0.620	0.702	-13.2	98	0.00	15.08
116	1,3,5-trimethylbenzene	2.135	2.508	-17.5	103	0.00	14.99
117	tert-butylbenzene	1.759	2.047	-16.4	99	0.00	15.33
118	pentachloroethane	0.480	0.633	-31.9#	116	0.00	15.42
119	1,2,4-trimethylbenzene	2.144	2.451	-14.3	101	0.00	15.38
120	sec-butylbenzene	2.811	3.383	-20.3#	103	0.00	15.55
121	1,3-dichlorobenzene	1.377	1.578	-14.6	100	0.00	15.73
122	p-isopropyltoluene	2.455	2.846	-15.9	102	0.00	15.66
123	1,4-dichlorobenzene	1.477	1.608	-8.9	100	0.00	15.81
124	1,2-dichlorobenzene	1.484	1.694	-14.2	101	0.00	16.19
125	1,4-Diethylbenzene			-----NA-----			
126	n-butylbenzene	1.285	1.444	-12.4	101	0.00	16.07
127	1,2,4,5-Tetramethylbenzen			-----NA-----			
128	1,2-dibromo-3-chloropropa	0.217	0.224	-3.2	97	0.00	16.92
129	1,3,5-trichlorobenzene	1.576	1.815	-15.2	100	0.00	17.09
130	1,2,4-trichlorobenzene	1.620	1.882	-16.2	100	0.00	17.66
131	hexachlorobutadiene	0.802	0.911	-13.6	107	0.00	17.76
132	naphthalene	3.432	3.909	-13.9	103	0.00	17.90
133	1,2,3-trichlorobenzene	1.682	1.980	-17.7	104	0.00	18.11
134	hexachloroethane	0.507	0.612	-20.7#	106	0.00	16.44
135	Benzyl chloride	1.442	1.320	8.5	90	0.00	15.92

(#) = Out of Range
3D112242.D M3D4810.M

SPCC's out = 0 CCC's out = 0
Wed Aug 12 09:32:33 2015 T



Continuing Calibration Summary

Page 1 of 3

Job Number: JC1106

Sample: V3D4824-CC4810

Account: UTC United Technologies Corporation

Lab FileID: 3D112563.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Evaluate Continuing Calibration Report

Data File : C:\msdchem\1\DATA\3D112563.D

Vial: 2

Acq On : 12 Aug 2015 9:17 am

Operator: ximenac

Sample : cc4810-20

Inst : MS3D

Misc : MS89457,V3D4824,5,,,,1

Multiplr: 1.00

MS Integration Params: rteint.p

Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)

Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

Last Update : Thu Aug 06 11:10:54 2015

Response via : Multiple Level Calibration

Min. RRF : 0.010 Min. Rel. Area : 50% Max. R.T. Dev 0.30min

Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)	R.T.
1 I	Tert Butyl Alcohol-d9	1.000	1.000	0.0	102	0.00	7.17
2	1,4-dioxane	0.081	0.102	-25.9#	118	0.00	11.04
3 M	tertiary butyl alcohol	1.080	1.037	4.0	95	-0.01	7.29
4 I	pentafluorobenzene	1.000	1.000	0.0	99	0.01	9.40
5	FREON 115			-----NA-----			
6	FREON 152A			-----NA-----			
7	chlorotrifluoroethene			-----NA-----			
8	chlorodifluoromethane	0.679	0.703	-3.5	102	0.00	3.86
9	dichlorodifluoromethane	0.909	1.008	-10.9	106	0.00	3.85
10	chloromethane	0.723	0.723	0.0	111	0.00	4.17
11	vinyl chloride	0.725	0.802	-10.6	111	0.00	4.41
12	bromomethane	0.468	0.527	-12.6	116	0.01	5.06
13	chloroethane	0.307	0.340	-10.7	117	0.02	5.23
14	1,3-butadiene			-----NA-----			
15	vinyl bromide			-----NA-----			
16	trichlorofluoromethane	0.756	0.923	-22.1#	118	0.01	5.68
17	pentane			-----NA-----			
18	ethyl ether	0.187	0.194	-3.7	103	0.00	6.07
19	freon 123a			-----NA-----			
20	FREON 123			-----NA-----			
21	acrolein	0.075	0.068	9.3	92	0.00	6.33
22	1,1-dichloroethene	0.595	0.594	0.2	97	0.00	6.49
23	acetone	0.112	0.123	-9.8	102	0.00	6.55
24	allyl chloride	0.185	0.187	-1.1	92	0.00	7.02
25	acetonitrile	0.037	0.039	-5.4	104	0.00	6.98
26 m	acetaldehyde			-----NA-----			
27	iodomethane	0.846	0.877	-3.7	96	0.00	6.77
28	iso-butyl alcohol	0.012	0.013	-8.3	102	0.01	9.69
29	carbon disulfide	1.504	1.576	-4.8	100	0.00	6.90
30	methylene chloride	0.446	0.459	-2.9	100	0.00	7.21
31	methyl acetate	0.249	0.263	-5.6	104	0.00	7.01
32	methyl tert butyl ether	1.320	1.233	6.6	90	0.00	7.54
33	trans-1,2-dichloroethene	0.543	0.565	-4.1	99	0.00	7.58
34	di-isopropyl ether	1.396	1.219	12.7	90	0.00	8.13
35	ethyl tert-butyl ether	1.315	1.195	9.1	88	0.01	8.59
36	2-butanone	0.036	0.034	5.6	87	0.01	8.86
37 M	1,1-dichloroethane	0.645	0.702	-8.8	100	0.01	8.15
38	chloroprene	0.472	0.428	9.3	92	0.01	8.26
39	acrylonitrile	0.129	0.142	-10.1	98	0.00	7.53
40	vinyl acetate	0.048	0.050	-4.2	101	0.00	8.14
41	ethyl acetate	0.049	0.040	18.4	77	0.01	8.87

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976

Continuing Calibration Summary

Page 2 of 3

Job Number: JC1106

Sample: V3D4824-CC4810

Account: UTC United Technologies Corporation

Lab FileID: 3D112563.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

42	2,2-dichloropropane	0.675	0.735	-8.9	111	0.00	8.88
43	cis-1,2-dichloroethene	0.436	0.414	5.0	96	0.01	8.89
44	propionitrile	0.048	0.052	-8.3	102	0.00	8.95
45	methyl acrylate	0.324	0.334	-3.1	92	0.01	8.95
46	bromochloromethane	0.210	0.227	-8.1	100	0.00	9.19
47	tetrahydrofuran	0.126	0.118	6.3	90	0.00	9.23
48	chloroform	0.622	0.652	-4.8	102	0.00	9.25
49	Tert-Butyl Formate	0.362	0.320	11.6	88	0.00	9.28
50 S	dibromofluoromethane (s)	0.383	0.404	-5.5	105	0.00	9.45
51 S	1,2-dichloroethane-d4 (s)	0.355	0.375	-5.6	106	0.00	9.87
52	freon 113	0.366	0.398	-8.7	109	0.02	6.47
53	methacrylonitrile	0.188	0.175	6.9	91	0.01	9.13
54	1,1,1-trichloroethane	0.642	0.722	-12.5	108	0.01	9.51
55	tert amyl alcohol			-----NA-----			
56	2,2,4-Trimethylpentane			-----NA-----			
57	tert-amyl methyl ether	1.314	1.165	11.3	92	0.00	9.97
58 I	1,4-difluorobenzene	1.000	1.000	0.0	98	0.01	10.32
59	epichlorohydrin	0.027	0.027	0.0	101	0.00	11.58
60	n-butyl alcohol	0.008	0.008#	0.0	102	0.02	10.46
61 M	cyclohexane	0.586	0.630	-7.5	107	0.00	9.57
62	carbon tetrachloride	0.484	0.591	-22.1#	112	0.00	9.71
63	1,1-dichloropropene	0.360	0.398	-10.6	100	0.00	9.68
64	hexane	0.322	0.284	11.8	91	0.00	7.88
65 M	benzene	1.175	1.258	-7.1	102	0.00	9.94
66	heptane	0.183	0.175	4.4	96	0.00	10.11
67	isopropyl acetate	0.511	0.505	1.2	94	0.01	9.87
68	1,2-dichloroethane	0.351	0.384	-9.4	100	0.00	9.96
69	trichloroethene	0.284	0.305	-7.4	103	0.01	10.66
70	ethyl acrylate			-----NA-----			
71	tert amyl ethyl ether			-----NA-----			
72	2-nitropropane	0.104	0.102	1.9	88	0.02	11.45
73	2-chloroethyl vinyl ether	0.180	0.181	-0.6	93	0.00	11.46
74	methyl methacrylate	0.402	0.426	-6.0	95	0.00	10.93
75	1,2-dichloropropane	0.307	0.334	-8.8	104	0.01	10.93
76	methylcyclohexane	0.531	0.563	-6.0	100	0.01	10.89
77	dibromomethane	0.181	0.209	-15.5	103	0.00	11.08
78	bromodichloromethane	0.374	0.406	-8.6	101	0.00	11.22
79	cis-1,3-dichloropropene	0.439	0.451	-2.7	91	0.00	11.68
80 S	toluene-d8 (s)	1.095	1.126	-2.8	100	0.00	11.97
81	4-methyl-2-pentanone	0.112	0.112	0.0	92	0.00	11.77
82	toluene	0.646	0.709	-9.8	97	0.01	12.04
83	3-methyl-1-butanol	0.008	0.008#	0.0	96	0.00	11.79
84	trans-1,3-dichloropropene	0.390	0.421	-7.9	95	0.00	12.24
85	ethyl methacrylate	0.345	0.292	15.4	78	0.01	12.24
86	1,1,2-trichloroethane	0.217	0.234	-7.8	100	0.00	12.45
87	2-hexanone	0.091	0.078	14.3	76	0.01	12.63
88 I	chlorobenzene-d5	1.000	1.000	0.0	99	0.00	13.48
89	tetrachloroethene	0.426	0.427	-0.2	91	0.00	12.63
90	1,3-dichloropropane	0.485	0.505	-4.1	96	0.01	12.64
91	butyl acetate	0.201	0.179	10.9	88	0.00	12.70
92	3,3-Dimethyl-1-Butanol	0.048	0.035	27.1#	84	0.00	12.79
93	dibromochloromethane	0.378	0.420	-11.1	99	0.00	12.90
94	1,2-dibromoethane	0.319	0.343	-7.5	94	0.00	13.05
95	n-Butyl Ether	1.513	1.253	17.2	79	0.00	13.43
96	chlorobenzene	0.897	0.973	-8.5	98	0.00	13.51
97	1,1,1-tetrachloroethane	0.398	0.445	-11.8	104	0.00	13.57
98	ethylbenzene	1.536	1.549	-0.8	95	0.00	13.56
99	m,p-xylene	0.574	0.605	-5.4	95	0.00	13.67

6.7.3
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Continuing Calibration Summary

Page 3 of 3

Job Number: JC1106

Sample: V3D4824-CC4810

Account: UTC United Technologies Corporation

Lab FileID: 3D112563.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

100	o-xylene	0.587	0.603	-2.7	91	0.00	14.09
101	styrene	0.987	0.913	7.5	85	0.01	14.10
102	Butyl Acrylate			-----NA-----			
103	bromoform	0.295	0.313	-6.1	95	0.00	14.36
104 I	1,4-dichlorobenzene-d4	1.000	1.000	0.0	90	0.00	15.79
105	isopropylbenzene	2.509	2.658	-5.9	90	0.00	14.43
106 S	4-bromofluorobenzene (s)	0.769	0.780	-1.4	93	0.00	14.63
107	bromobenzene	0.741	0.806	-8.8	95	0.01	14.83
		----- True	Calc.	% Drift			
108	cyclohexanone	200.000	113.675	43.2#	24	0.01	14.60
		----- AvgRF	CCRF	% Dev			
109	1,1,2,2-tetrachloroethane	0.726	0.826	-13.8	98	0.00	14.73
110	trans-1,4-dichloro-2-bute	0.161	0.134	16.8	71	0.01	14.78
111	1,2,3-trichloropropane	0.166	0.191	-15.1	100	0.00	14.81
112	n-propylbenzene	2.732	2.954	-8.1	91	0.00	14.84
113	4-Ethyltoluene			-----NA-----			
114	2-chlorotoluene	0.638	0.702	-10.0	95	0.00	14.99
115	4-chlorotoluene	0.620	0.667	-7.6	91	0.00	15.08
116	1,3,5-trimethylbenzene	2.135	2.337	-9.5	93	0.00	14.99
117	tert-butylbenzene	1.759	1.827	-3.9	88	0.00	15.34
118	pentachloroethane	0.480	0.600	-25.0#	112	0.00	15.42
119	1,2,4-trimethylbenzene	2.144	2.270	-5.9	90	0.00	15.38
120	sec-butylbenzene	2.811	3.089	-9.9	91	0.00	15.55
121	1,3-dichlorobenzene	1.377	1.505	-9.3	90	0.00	15.73
122	p-isopropyltoluene	2.455	2.623	-6.8	90	0.00	15.67
123	1,4-dichlorobenzene	1.477	1.534	-3.9	90	0.00	15.82
124	1,2-dichlorobenzene	1.484	1.598	-7.7	91	0.00	16.20
125	1,4-Diethylbenzene			-----NA-----			
126	n-butylbenzene	1.285	1.304	-1.5	86	0.00	16.07
127	1,2,4,5-Tetramethylbenzen			-----NA-----			
128	1,2-dibromo-3-chloropropa	0.217	0.208	4.1	82	0.00	16.93
129	1,3,5-trichlorobenzene	1.576	1.671	-6.0	87	0.00	17.10
130	1,2,4-trichlorobenzene	1.620	1.720	-6.2	89	0.00	17.66
131	hexachlorobutadiene	0.802	0.855	-6.6	95	0.00	17.76
132	naphthalene	3.432	3.601	-4.9	89	0.00	17.90
133	1,2,3-trichlorobenzene	1.682	1.860	-10.6	94	0.00	18.12
134	hexachloroethane	0.507	0.556	-9.7	101	0.00	16.44
135	Benzyl chloride	1.442	1.368	5.1	91	0.00	15.93

(#) = Out of Range
3D112241.D M3D4810.M

SPCC's out = 0 CCC's out = 0
Wed Aug 12 14:29:13 2015 3D



New Jersey
ACCUTEST[®]
LABORATORIES

GC/MS Volatiles

Raw Data

7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\

Data File : 3D112575.D

Acq On : 12 Aug 2015 2:48 pm

Operator : ximenac

Sample : jc1106-1

Misc : MS89468,V3D4824,5,,,1

ALS Vial : 14 Sample Multiplier: 1

Quant Time: Aug 13 10:35:00 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

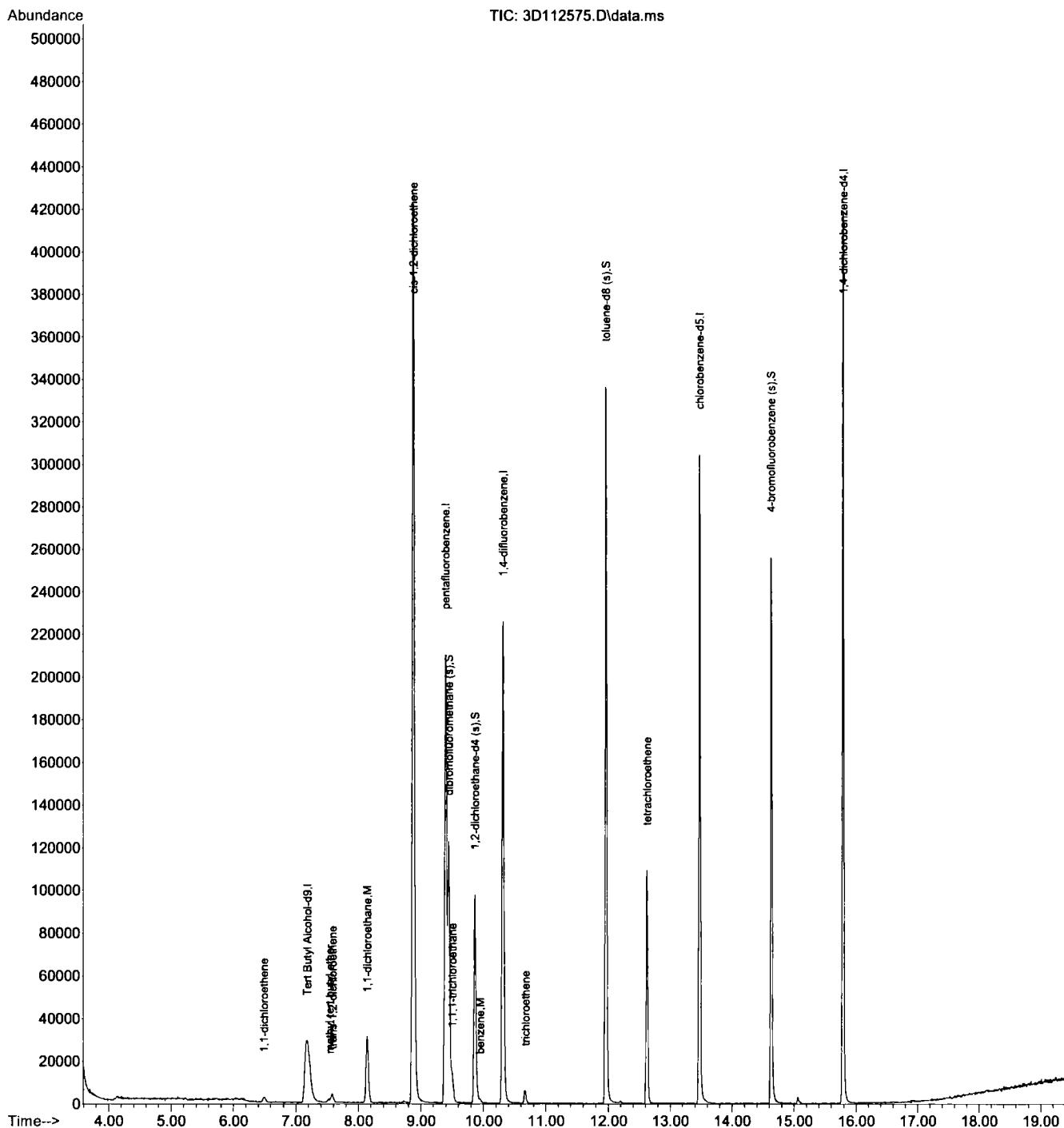
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.183	65	93640	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	196615	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	224320	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	180779	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	111990	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	84286	55.95	ug/L	0.00
Spiked Amount 50.000 Range	76 - 120		Recovery	=	111.90%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	77038	55.21	ug/L	0.00
Spiked Amount 50.000 Range	73 - 122		Recovery	=	110.42%	
80) toluene-d8 (s)	11.960	98	243477	49.56	ug/L	0.00
Spiked Amount 50.000 Range	84 - 119		Recovery	=	99.12%	
106) 4-bromofluorobenzene (s)	14.629	95	83190	48.28	ug/L	0.00
Spiked Amount 50.000 Range	78 - 117		Recovery	=	96.56%	
Target Compounds						
				Qvalue		
22) 1,1-dichloroethene	6.491	61	2534	1.08	ug/L	97
32) methyl tert butyl ether	7.540	73	3296	0.63	ug/L	97
33) trans-1,2-dichloroethene	7.587	61	2850	1.34	ug/L	94
37) 1,1-dichloroethane	8.143	63	41747	16.47	ug/L	98
43) cis-1,2-dichloroethene	8.877	96	239895	140.01	ug/L	88
54) 1,1,1-trichloroethane	9.501	97	10211	4.05	ug/L	97
65) benzene	9.947	78	1423	0.27	ug/L	81
69) trichloroethene	10.670	95	2628	2.06	ug/L	89
89) tetrachloroethene	12.626	166	37409	24.30	ug/L	98

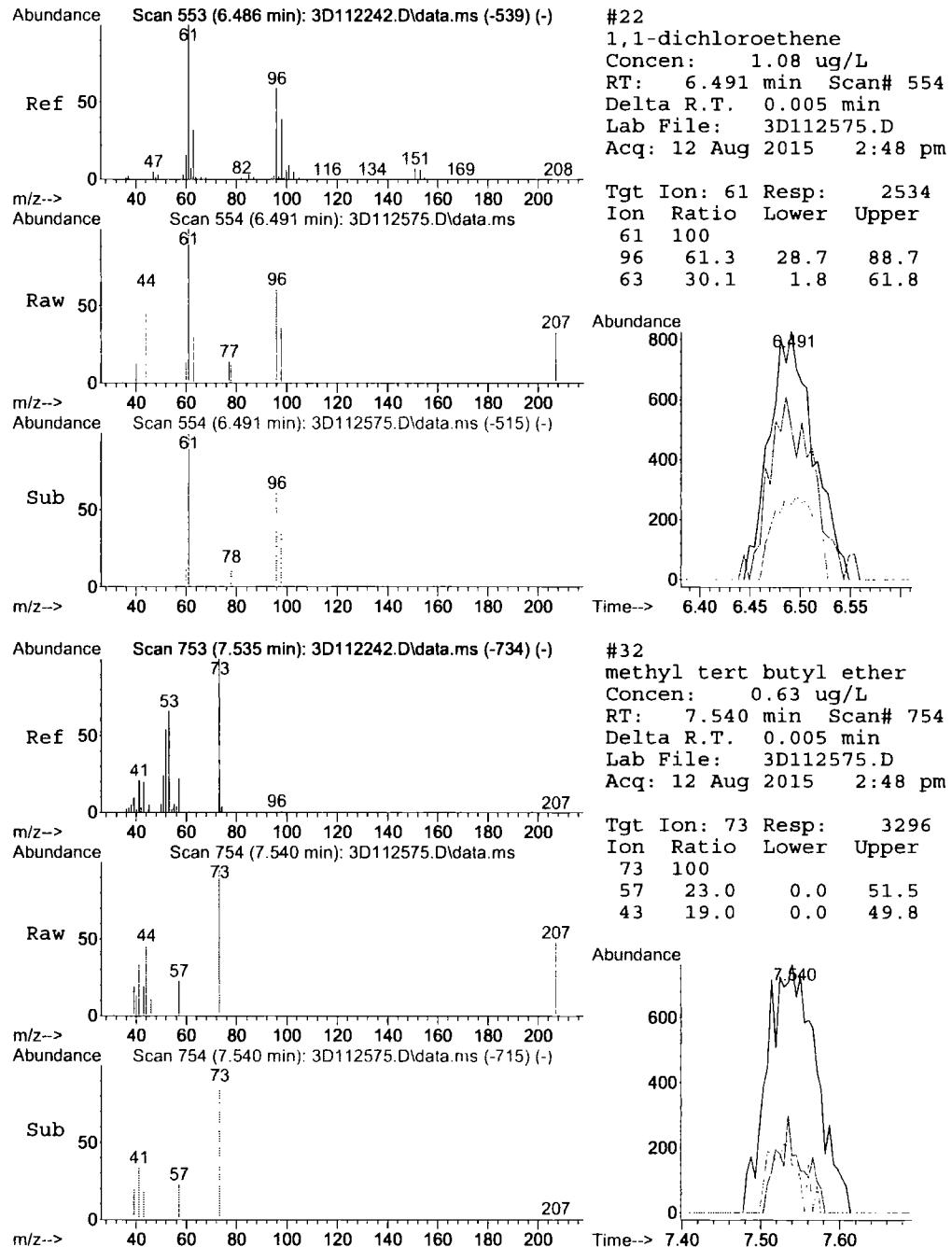
(#= qualifier out of range (m) = manual integration (+) = signals summed

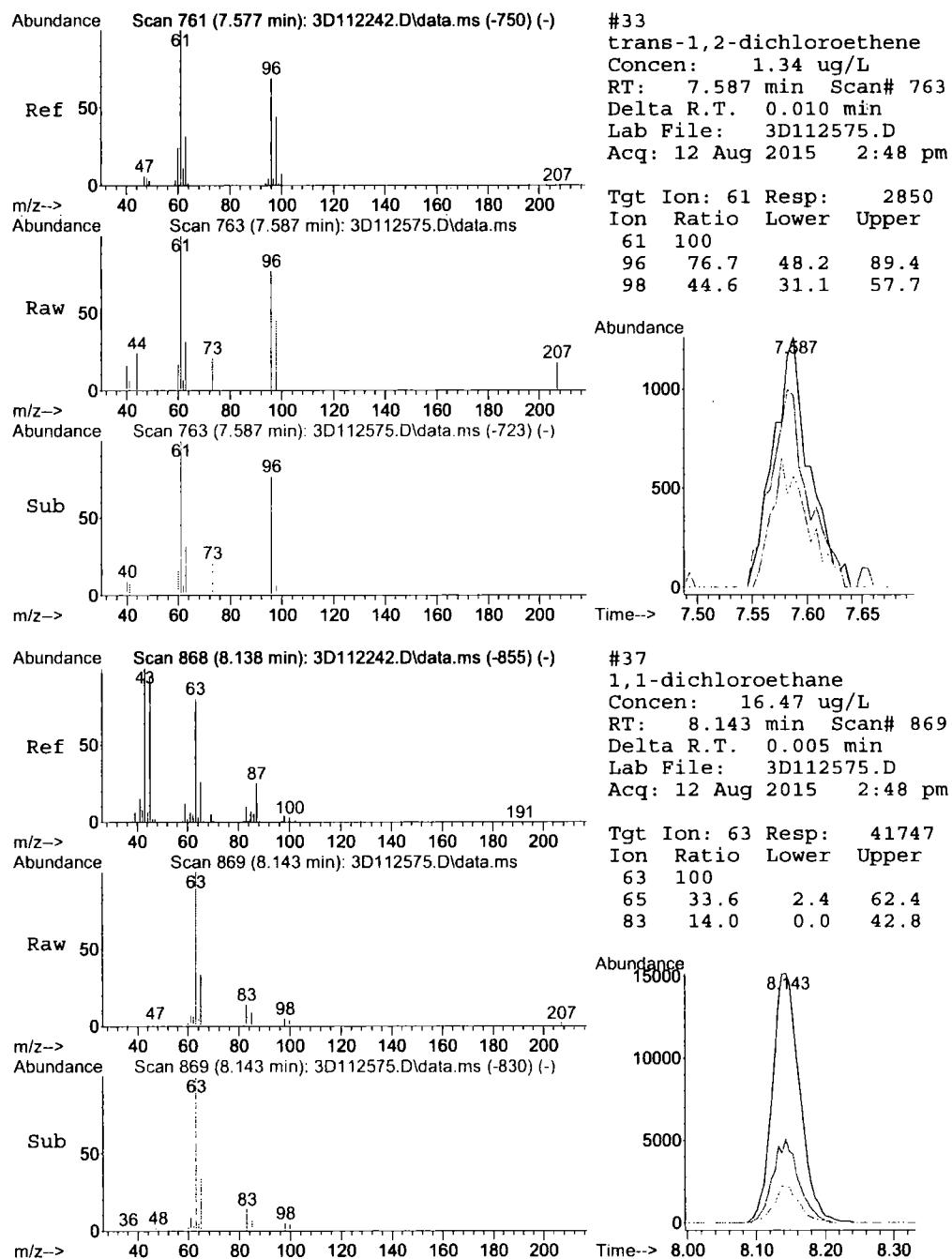
Quantitation Report (QT Reviewed)

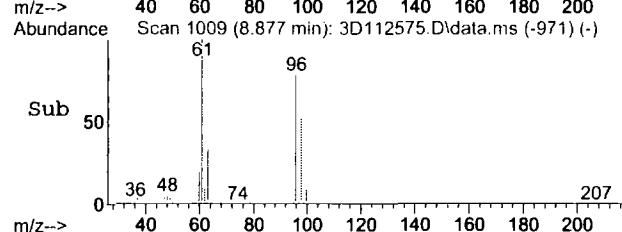
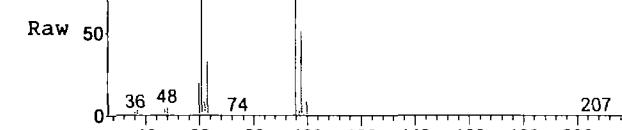
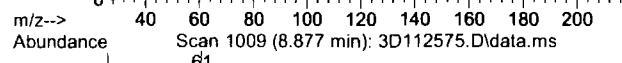
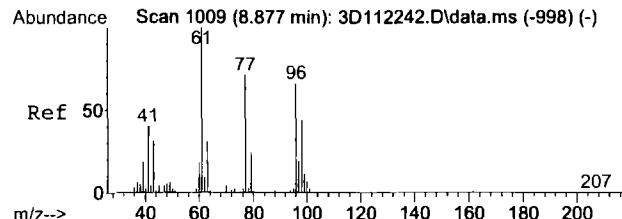
Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
 Data File : 3D112575.D
 Acq On : 12 Aug 2015 2:48 pm
 Operator : ximenac
 Sample : jc1106-1
 Misc : MS89468,V3D4824,5,,,1
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Aug 13 10:35:00 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



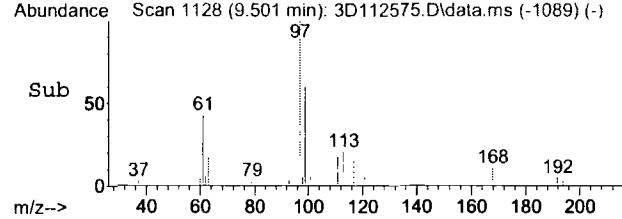
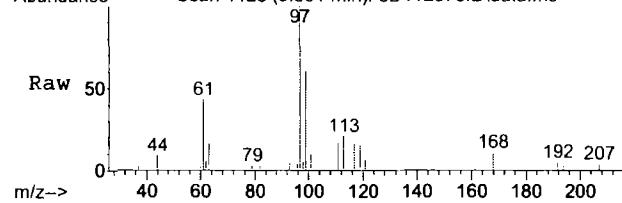
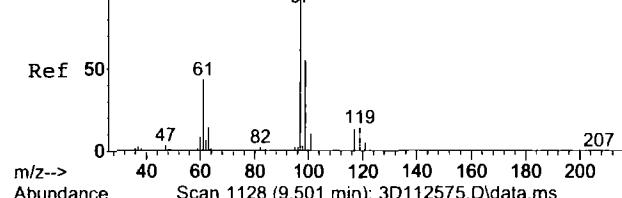
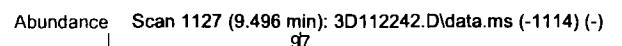
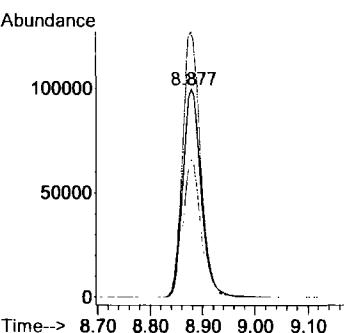






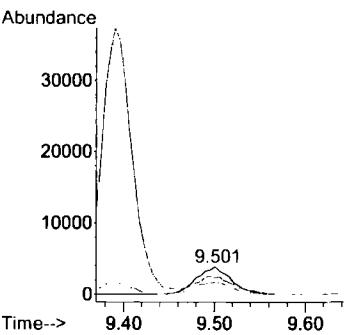
#43
cis-1,2-dichloroethene
Concen: 140.01 ug/L
RT: 8.877 min Scan# 1009
Delta R.T. -0.000 min
Lab File: 3D112575.D
Acq: 12 Aug 2015 2:48 pm

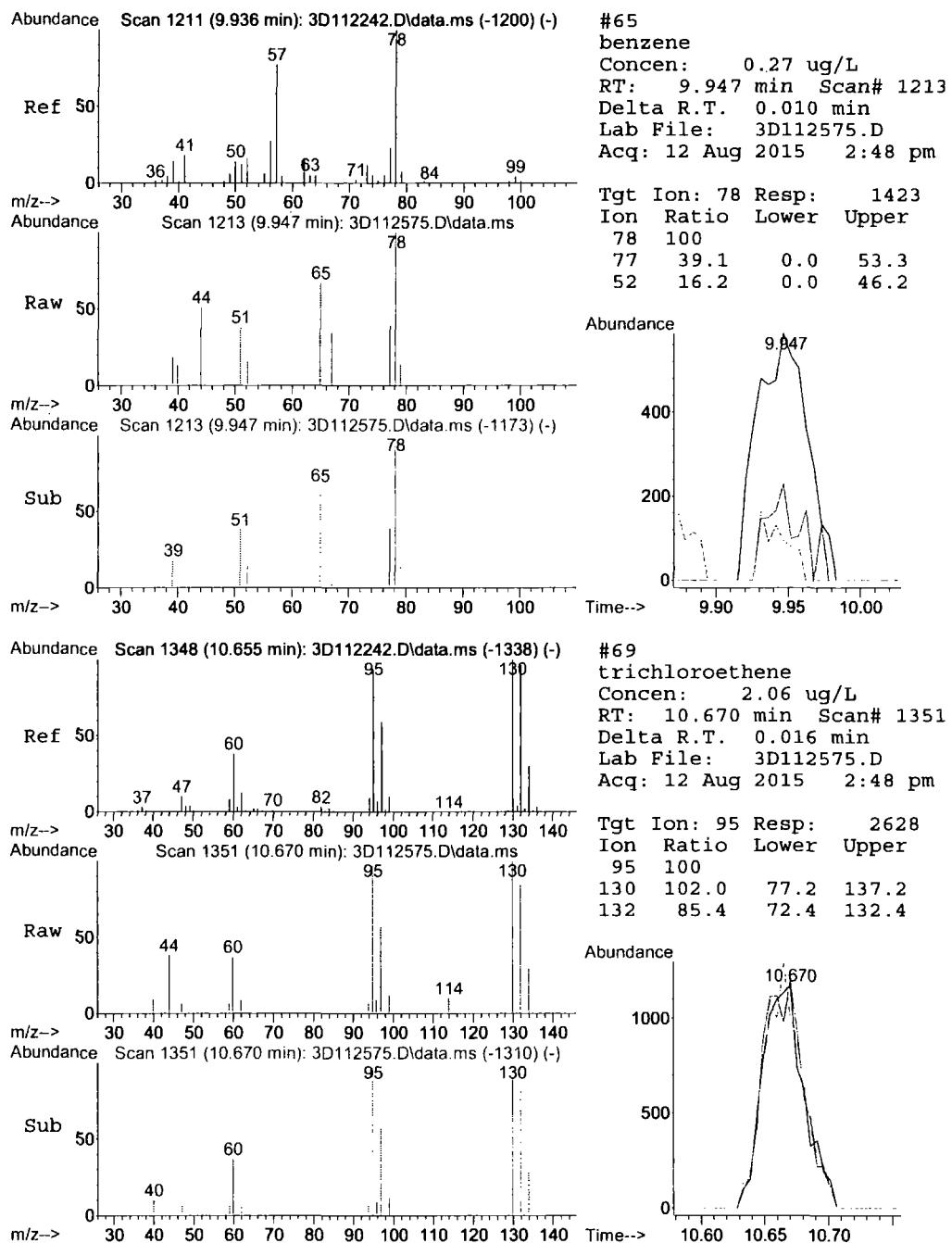
Tgt	Ion:	96	Resp:	239895
Ion	Ratio		Lower	Upper
96	100			
61	128.1		120.5	180.5
98	66.2		36.4	96.4

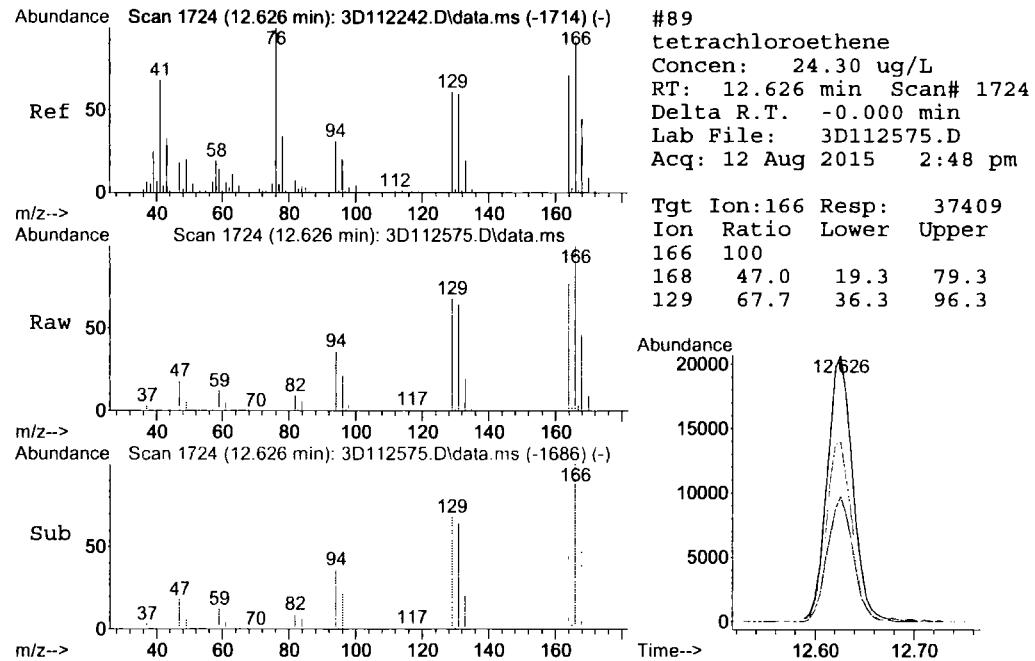


#54
1,1,1-trichloroethane
Concen: 4.05 ug/L
RT: 9.501 min Scan# 1128
Delta R.T. 0.005 min
Lab File: 3D112575.D
Acq: 12 Aug 2015 2:48 pm

Tgt	Ion:	97	Resp:	10211
Ion	Ratio		Lower	Upper
97	100			
99	59.5		33.7	93.7
61	43.0		13.5	73.5







Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
 Data File : 3D112574.D
 Acq On : 12 Aug 2015 2:21 pm
 Operator : ximenac
 Sample : jc1106-2
 Misc : MS89468,V3D4824,5,,,1
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 13 09:53:50 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.168	65	92619	500.00	ug/L	-0.01
4) pentafluorobenzene	9.391	168	180692	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	212538	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	175332	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	108820	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	77640	56.08	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	= 112.16%		
51) 1,2-dichloroethane-d4 (s)	9.863	65	73686	57.46	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	= 114.92%		
80) toluene-d8 (s)	11.960	98	231940	49.83	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	= 99.66%		
106) 4-bromofluorobenzene (s)	14.629	95	79559	47.52	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	= 95.04%		
Target Compounds						
32) methyl tert butyl ether	7.530	73	1010	0.21	ug/L	72
37) 1,1-dichloroethane	8.143	63	15416	6.62	ug/L	96
43) cis-1,2-dichloroethene	8.877	96	5772	3.67	ug/L	82
54) 1,1,1-trichloroethane	9.496	97	14745	6.36	ug/L	99
69) trichloroethene	10.660	95	2162	1.79	ug/L	79
89) tetrachloroethene	12.626	166	27341	18.32	ug/L	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\

Data File : 3D112574.D

Acq On : 12 Aug 2015 2:21 pm

Operator : ximenac

Sample : jc1106-2

Misc : MS89468,V3D4824,5,,,1

ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 13 09:53:50 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

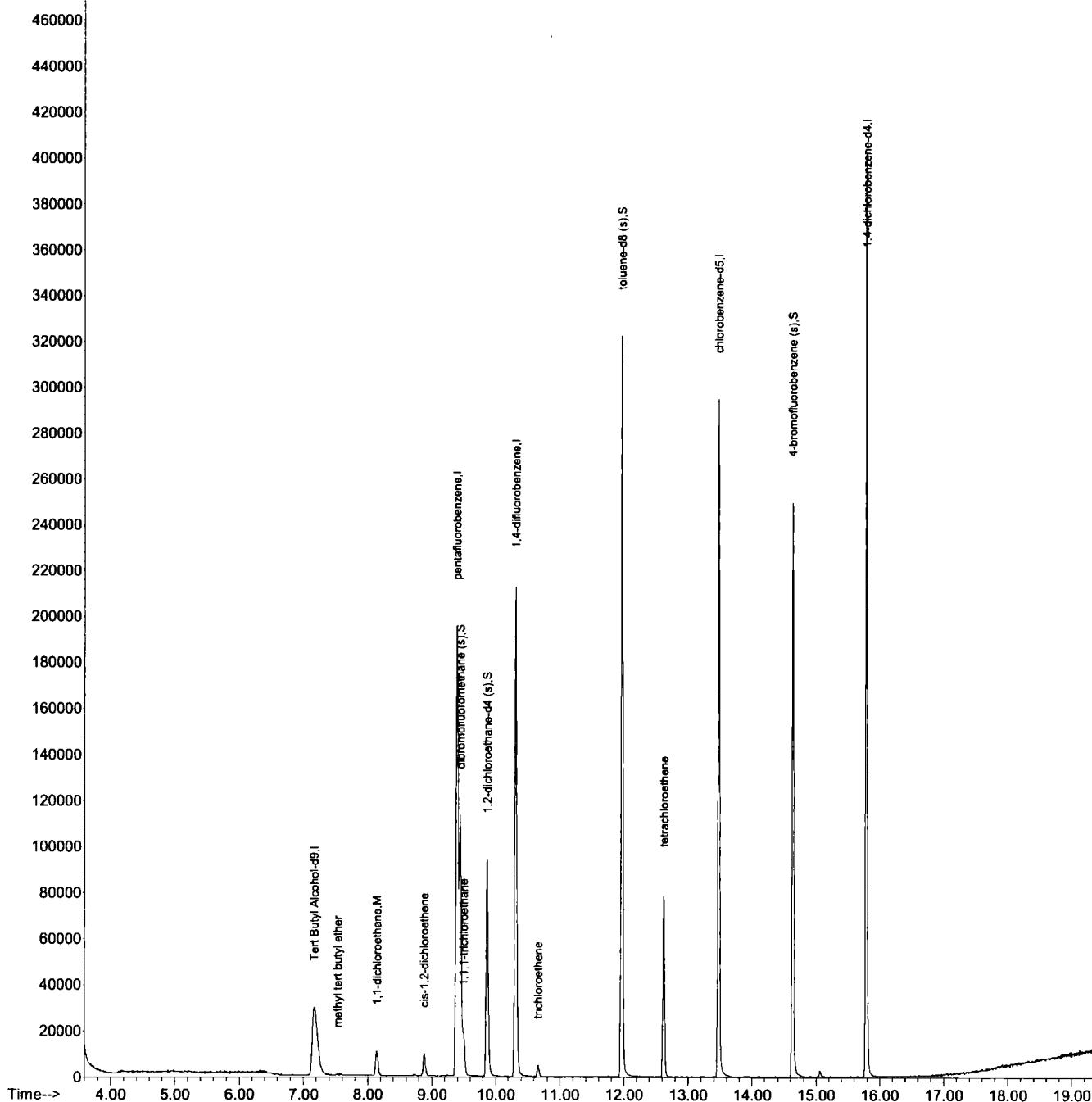
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

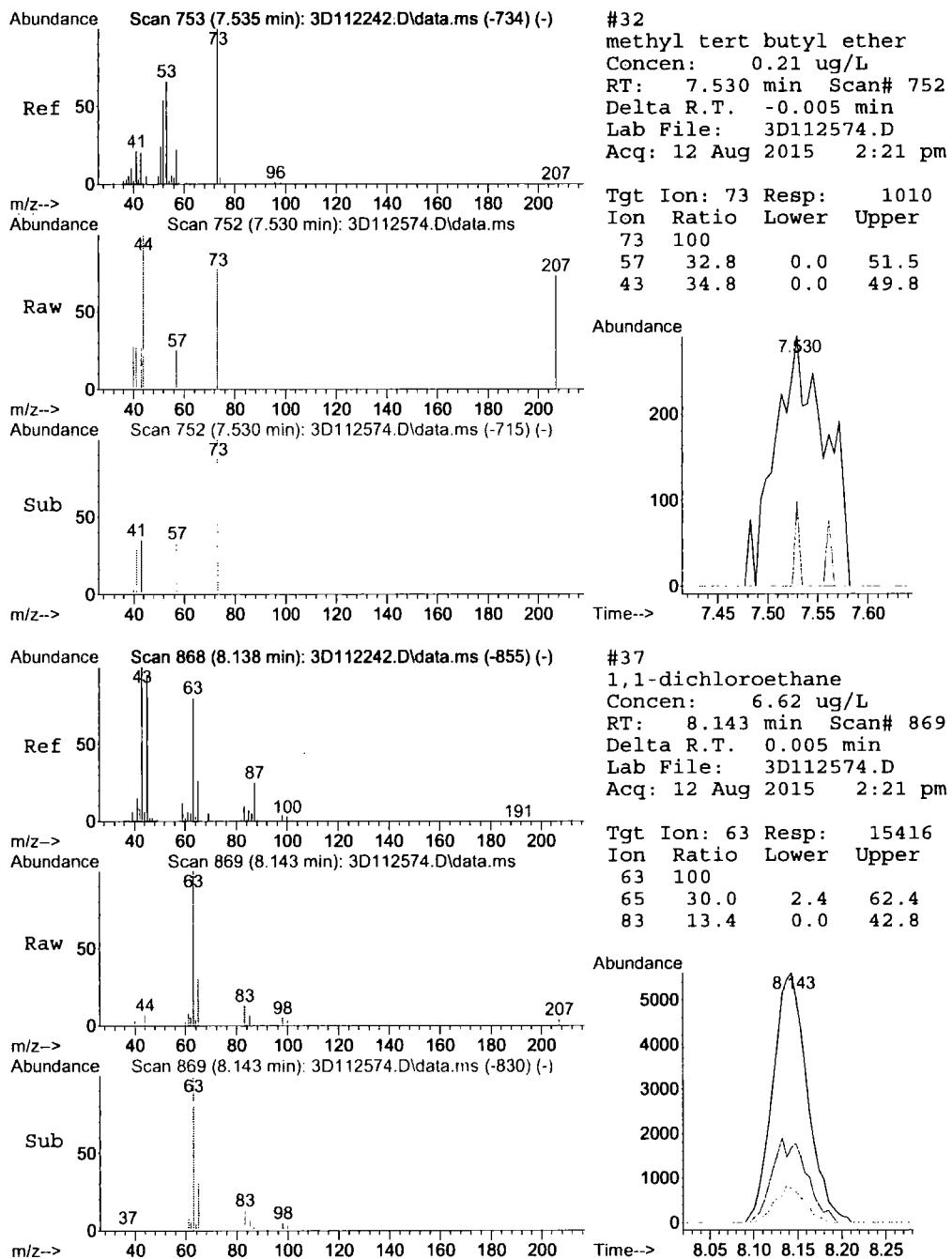
Abundance
480000

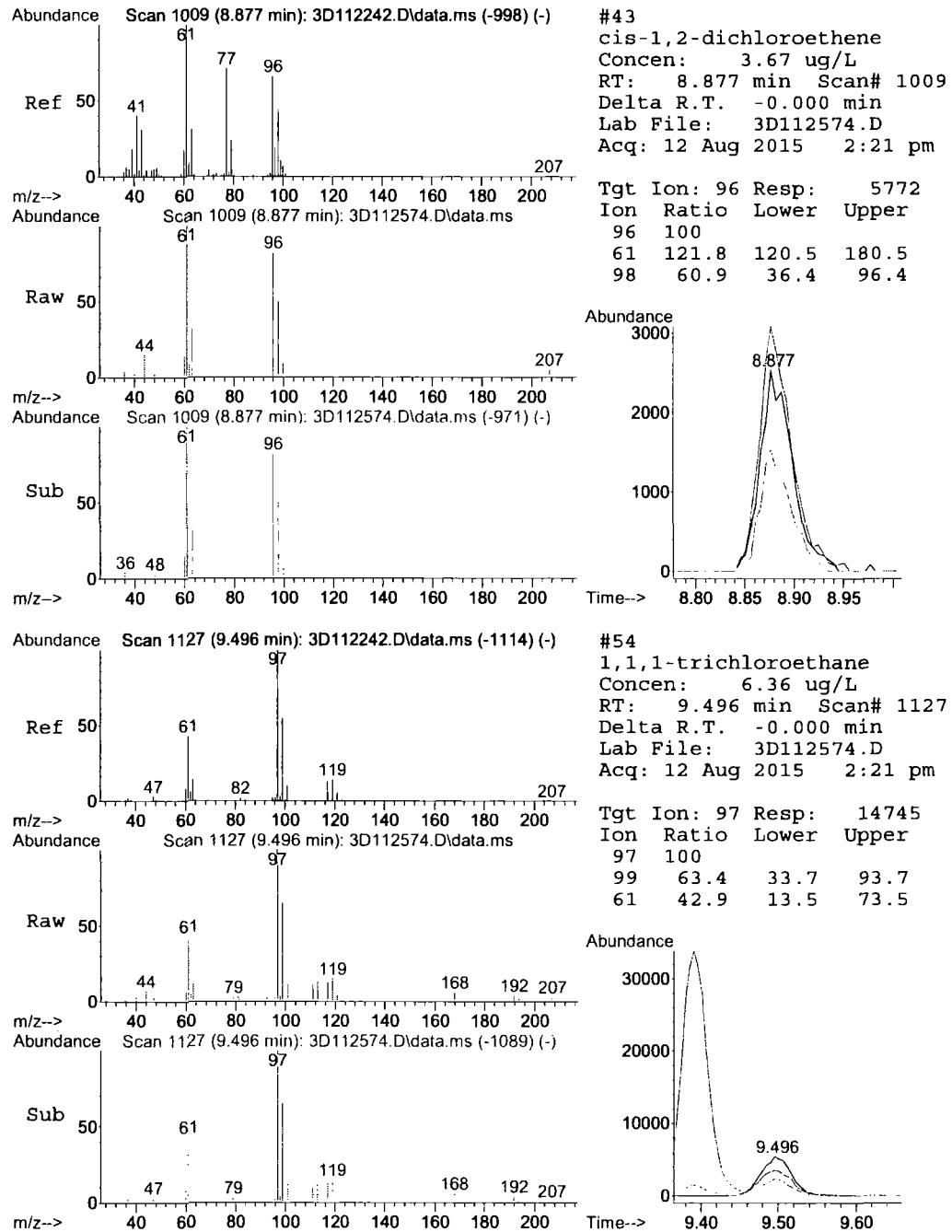
TIC: 3D112574.D\data.ms

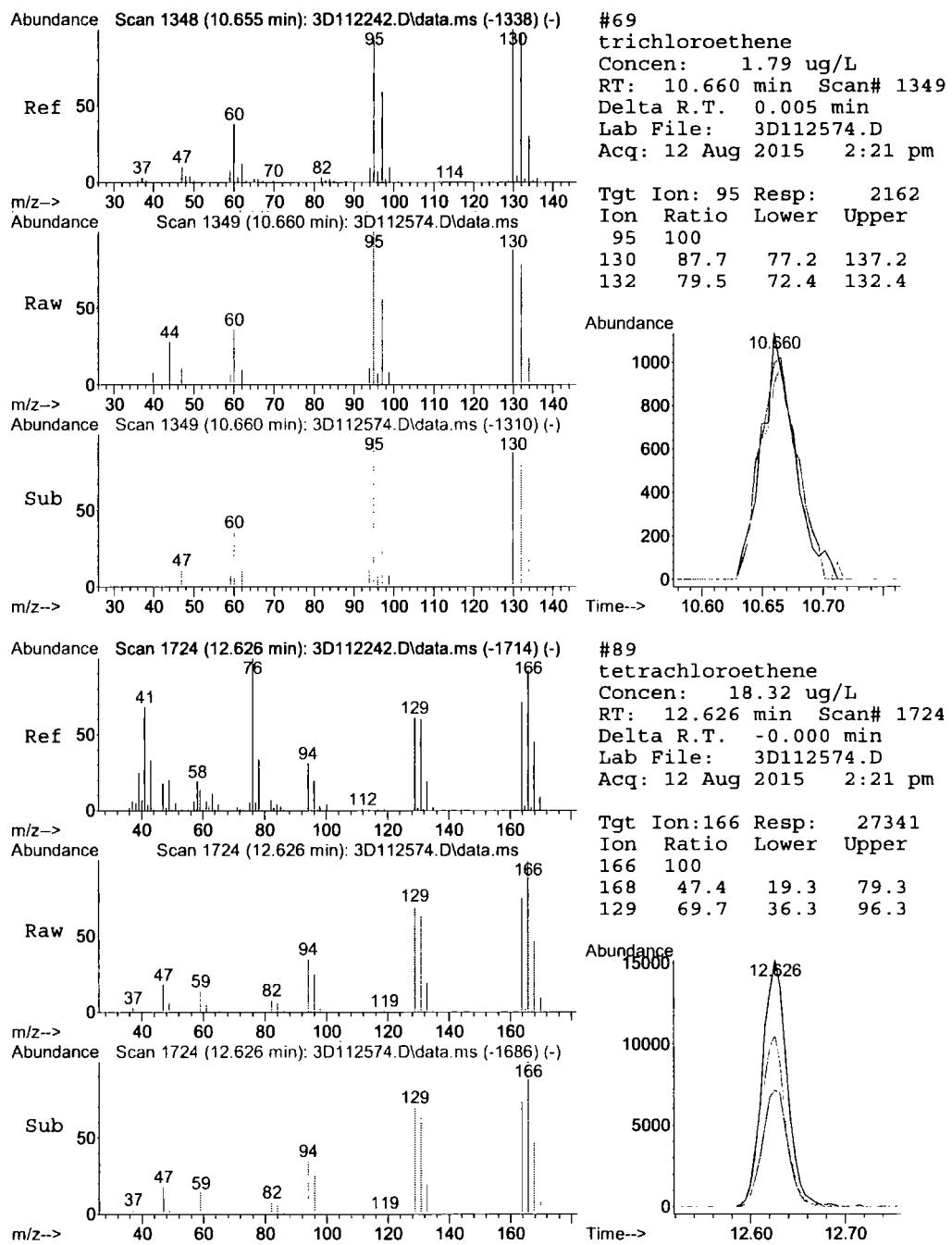


7.12

7







Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112531.D
 Acq On : 11 Aug 2015 6:18 pm
 Operator : ximenac
 Sample : jc1106-3
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Aug 12 09:28:45 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

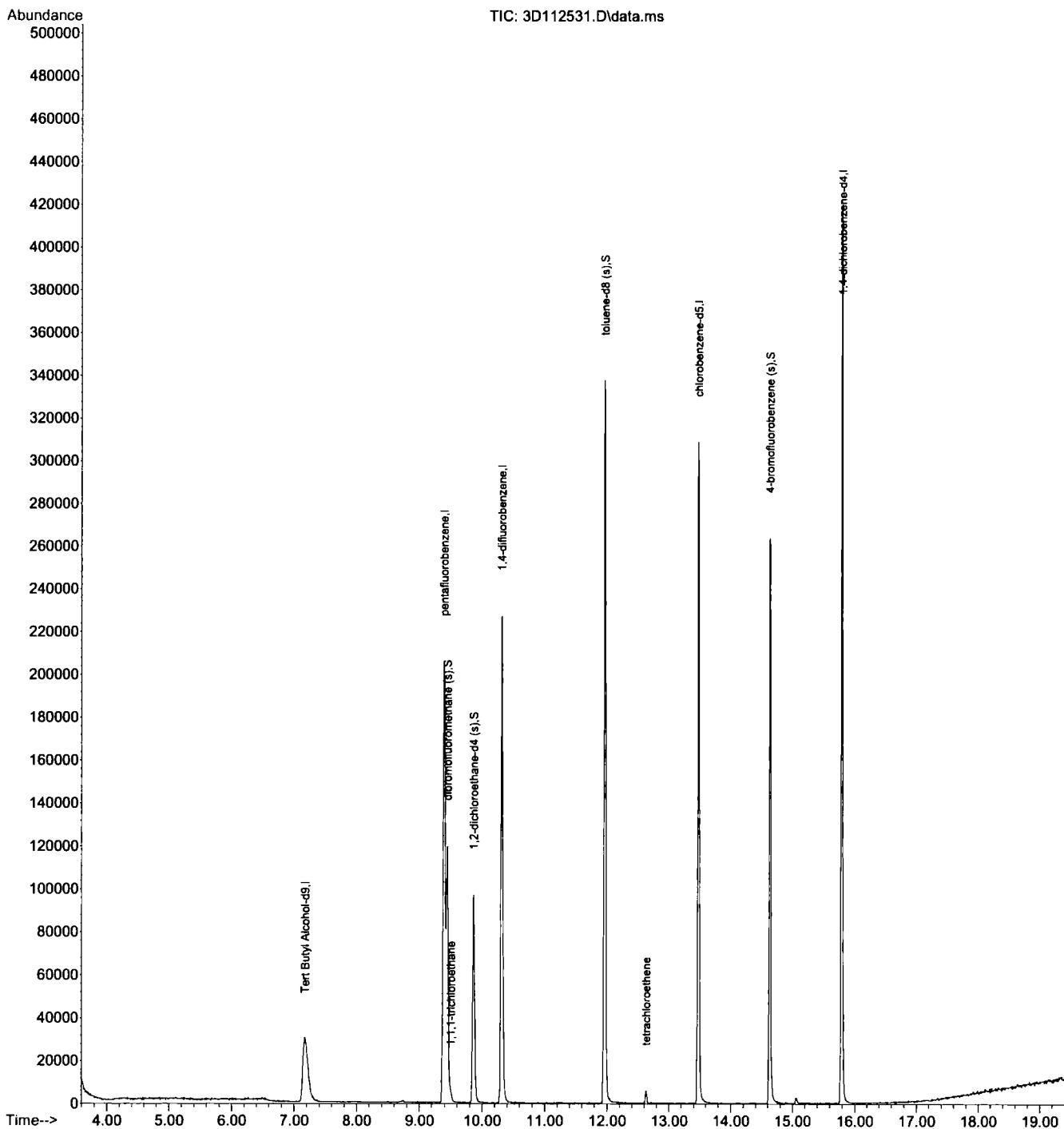
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	93947	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	191420	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	223101	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	183270	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	111855	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	81015	55.24	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	= 110.48%		
51) 1,2-dichloroethane-d4 (s)	9.863	65	77053	56.71	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	= 113.42%		
80) toluene-d8 (s)	11.960	98	246220	50.39	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	= 100.78%		
106) 4-bromofluorobenzene (s)	14.629	95	84059	48.84	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	= 97.68%		
Target Compounds						
				Qvalue		
54) 1,1,1-trichloroethane	9.501	97	3314	1.35	ug/L	84
89) tetrachloroethene	12.631	166	2096	1.34	ug/L	96

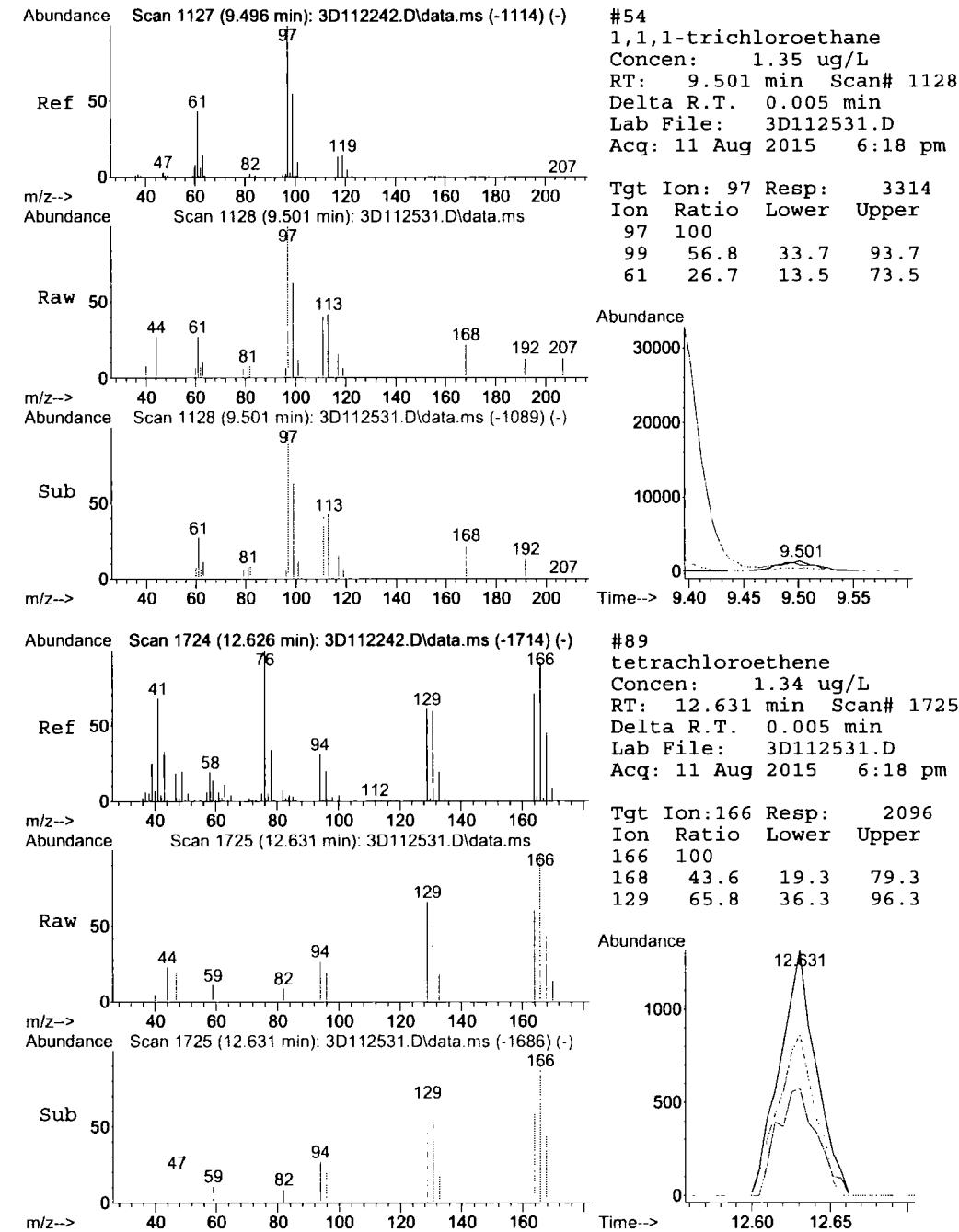
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112531.D
 Acq On : 11 Aug 2015 6:18 pm
 Operator : ximenac
 Sample : jc1106-3
 Misc : MS89468,V3D4822,5,,,,1
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Aug 12 09:28:45 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112532.D
 Acq On : 11 Aug 2015 6:45 pm
 Operator : ximenac
 Sample : jc1106-4
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Aug 12 09:29:07 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

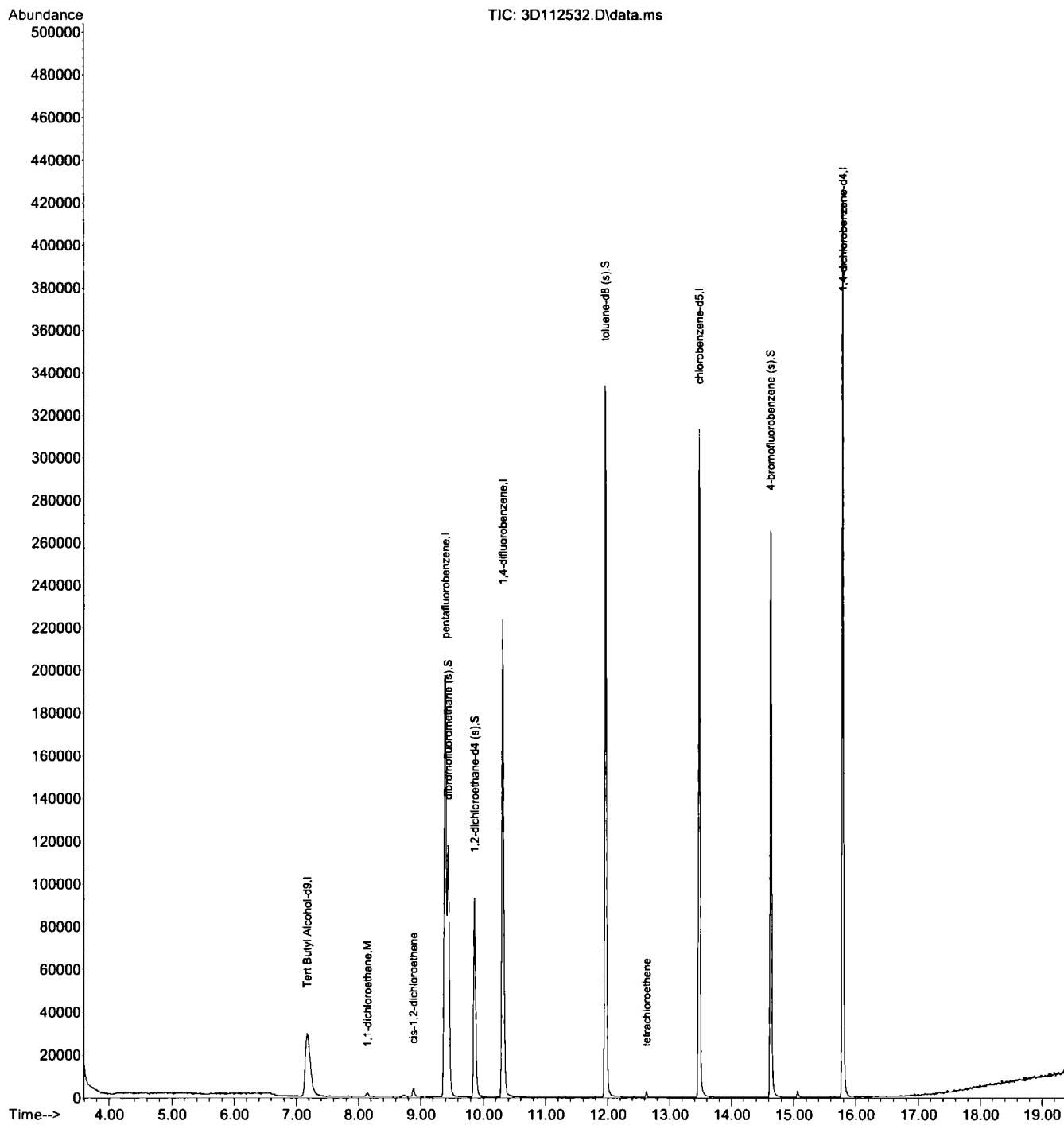
Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	93695	500.00	ug/L	0.00
4) pentafluorobenzene	9.396	168	187081	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	219575	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	184978	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	115133	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.438	113	81663	56.97	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	=	113.94%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	75123	56.58	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	=	113.16%	
80) toluene-d8 (s)	11.960	98	243928	50.72	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	=	101.44%	
106) 4-bromofluorobenzene (s)	14.629	95	85636	48.34	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	=	96.68%	
Target Compounds						
				Qvalue		
37) 1,1-dichloroethane	8.143	63	2696	1.12	ug/L	95
43) cis-1,2-dichloroethene	8.882	96	2277	1.40	ug/L	# 72
89) tetrachloroethene	12.631	166	961	0.61	ug/L	85

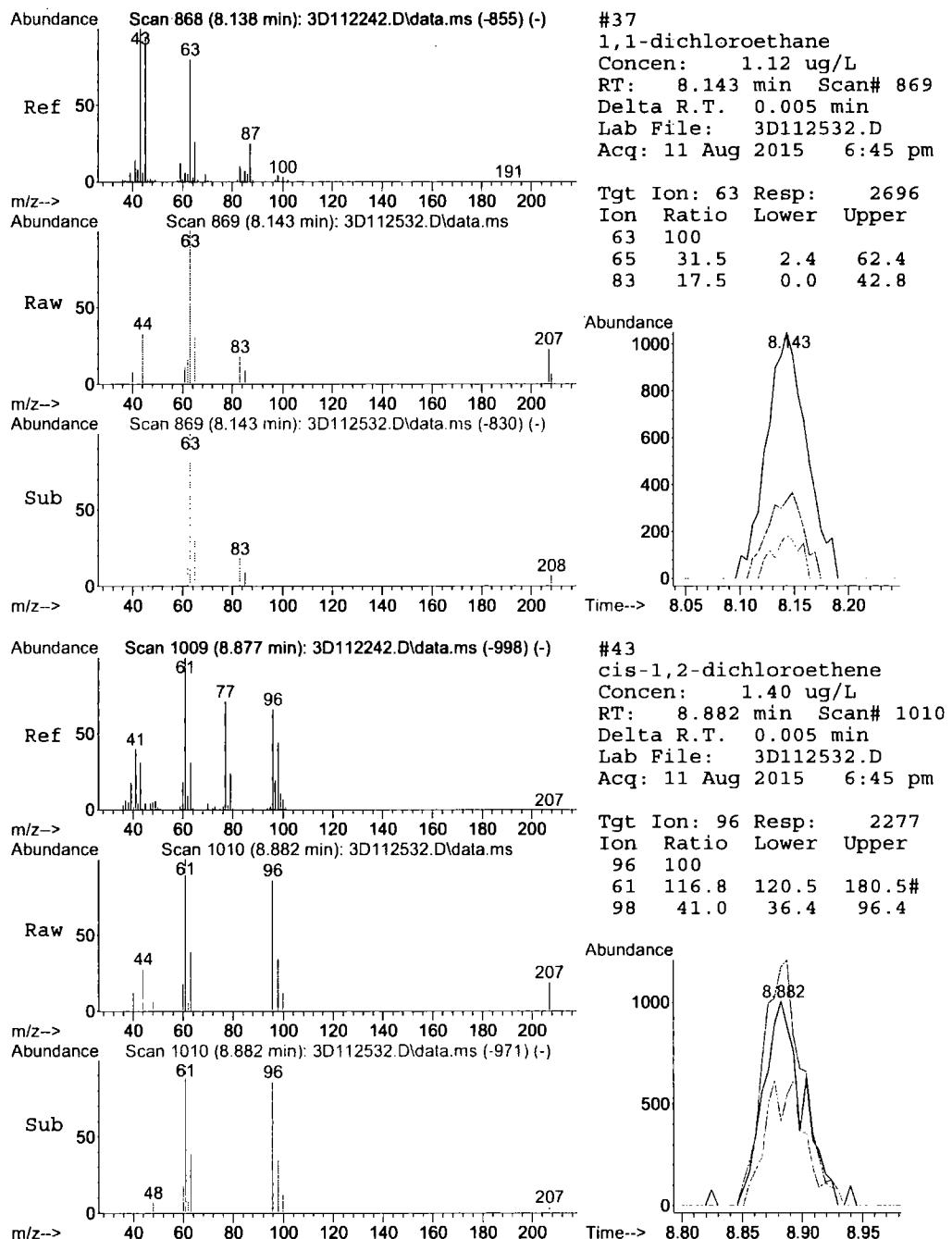
(#) = qualifier out of range (m) = manual integration (+) = signals summed

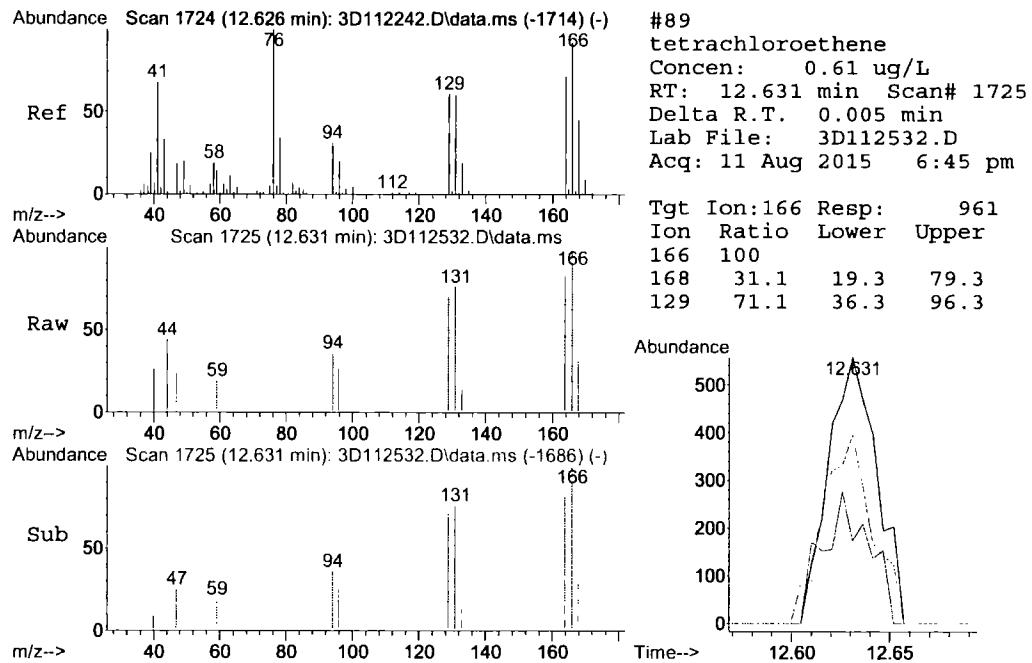
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112532.D
 Acq On : 11 Aug 2015 6:45 pm
 Operator : ximenac
 Sample : jc1106-4
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Aug 12 09:29:07 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration







Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112523.D
 Acq On : 11 Aug 2015 2:36 pm
 Operator : ximenac
 Sample : jc1106-5
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Aug 11 16:11:25 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.173	65	97975	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	195078	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	228938	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	186093	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	118621	50.00	ug/L	0.00

System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	83641	55.96	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	111.92%
51) 1,2-dichloroethane-d4 (s)	9.863	65	78760	56.88	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	113.76%
80) toluene-d8 (s)	11.960	98	252606	50.38	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	100.76%
106) 4-bromofluorobenzene (s)	14.629	95	88742	48.62	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.24%

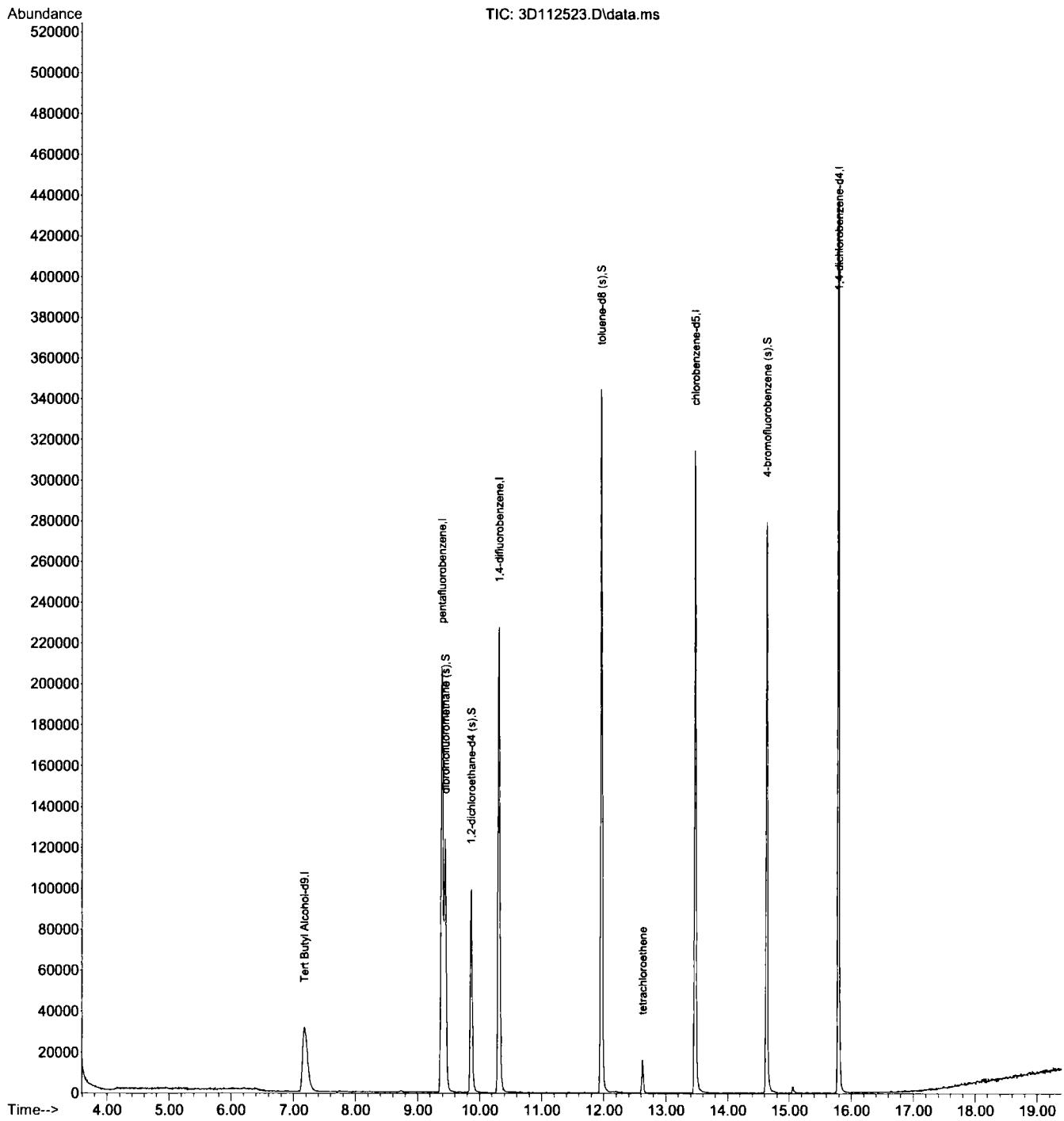
Target Compounds					Qvalue
89) tetrachloroethene	12.626	166	5806	3.66	ug/L

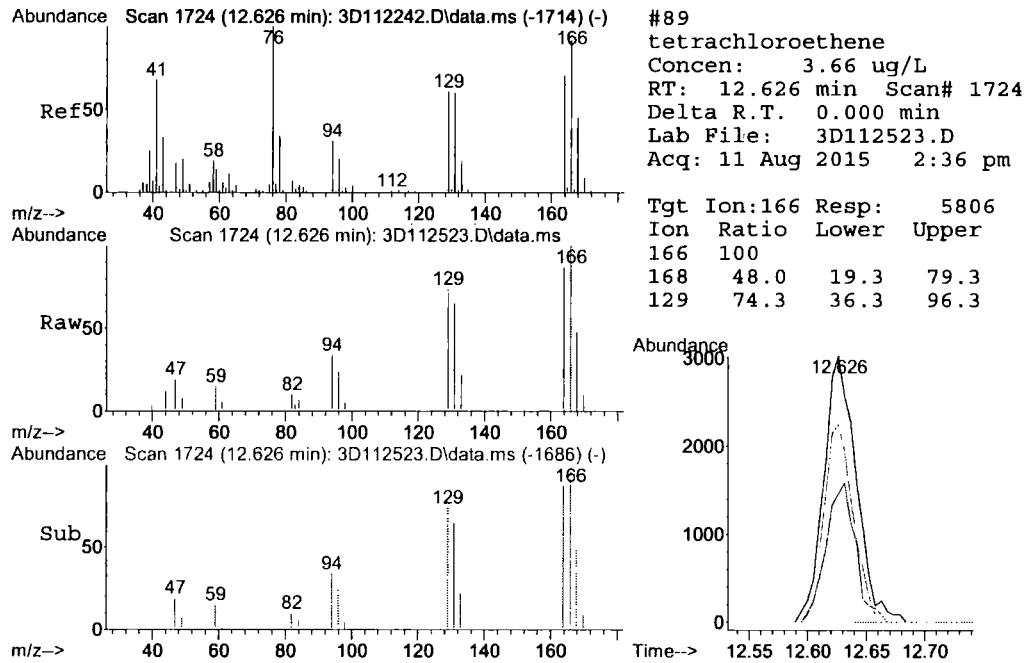
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112523.D
 Acq On : 11 Aug 2015 2:36 pm
 Operator : ximenac
 Sample : jc1106-5
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Aug 11 16:11:25 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\

Data File : 3D112530.D

Acq On : 11 Aug 2015 5:51 pm

Operator : ximenac

Sample : jc1106-6

Misc : MS89468,V3D4822,5,,,1

ALS Vial : 23 Sample Multiplier: 1

Quant Time: Aug 12 09:28:33 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

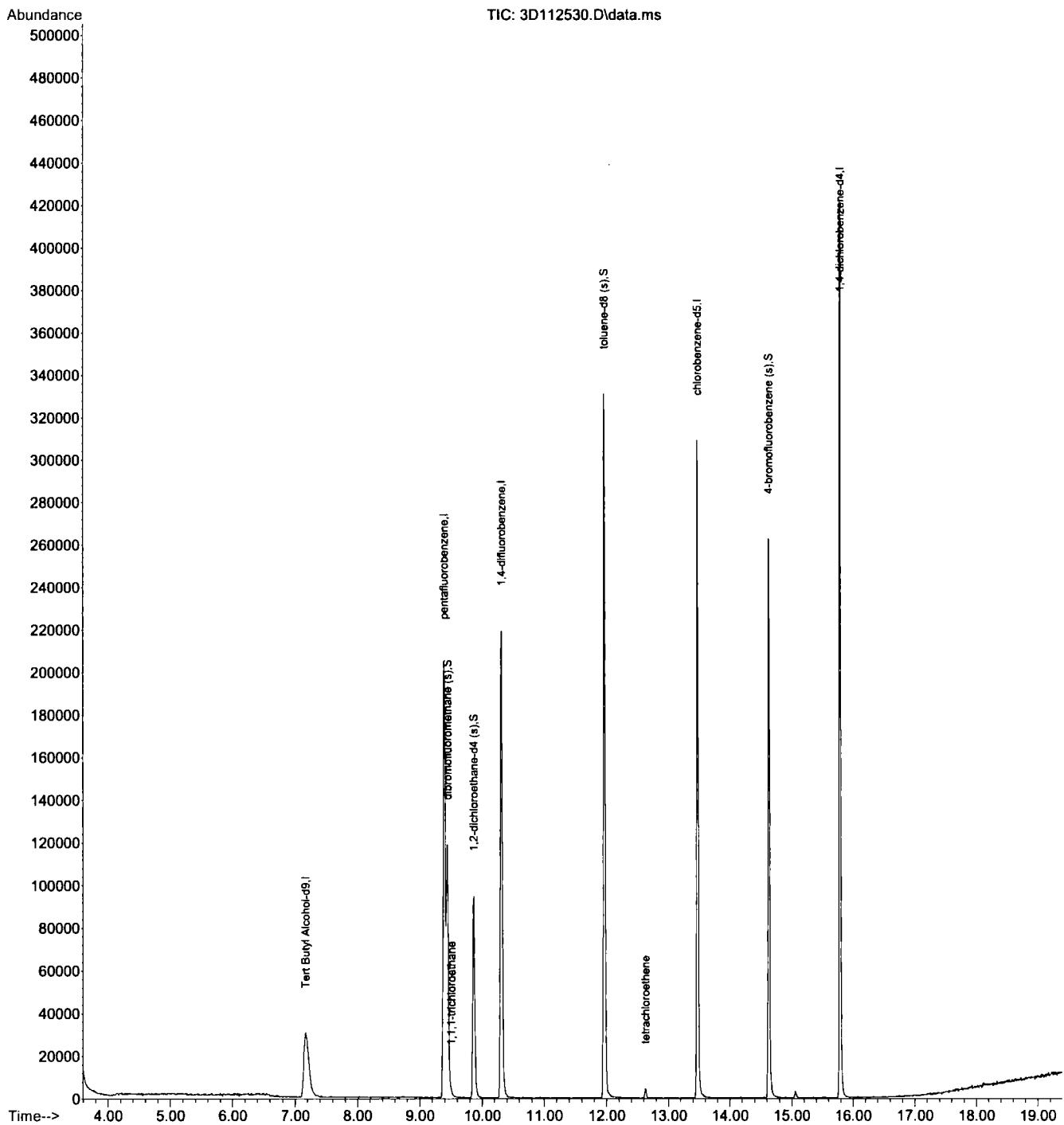
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.168	65	95232	500.00	ug/L	-0.01
4) pentafluorobenzene	9.391	168	188915	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	220873	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	182140	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	113662	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	80541	55.64	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	=	111.28%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	75278	56.14	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	=	112.28%	
80) toluene-d8 (s)	11.960	98	242380	50.11	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	=	100.22%	
106) 4-bromofluorobenzene (s)	14.629	95	83691	47.86	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	=	95.72%	
Target Compounds						
54) 1,1,1-trichloroethane	9.506	97	2611	1.08	ug/L	86
89) tetrachloroethene	12.626	166	1659	1.07	ug/L	88

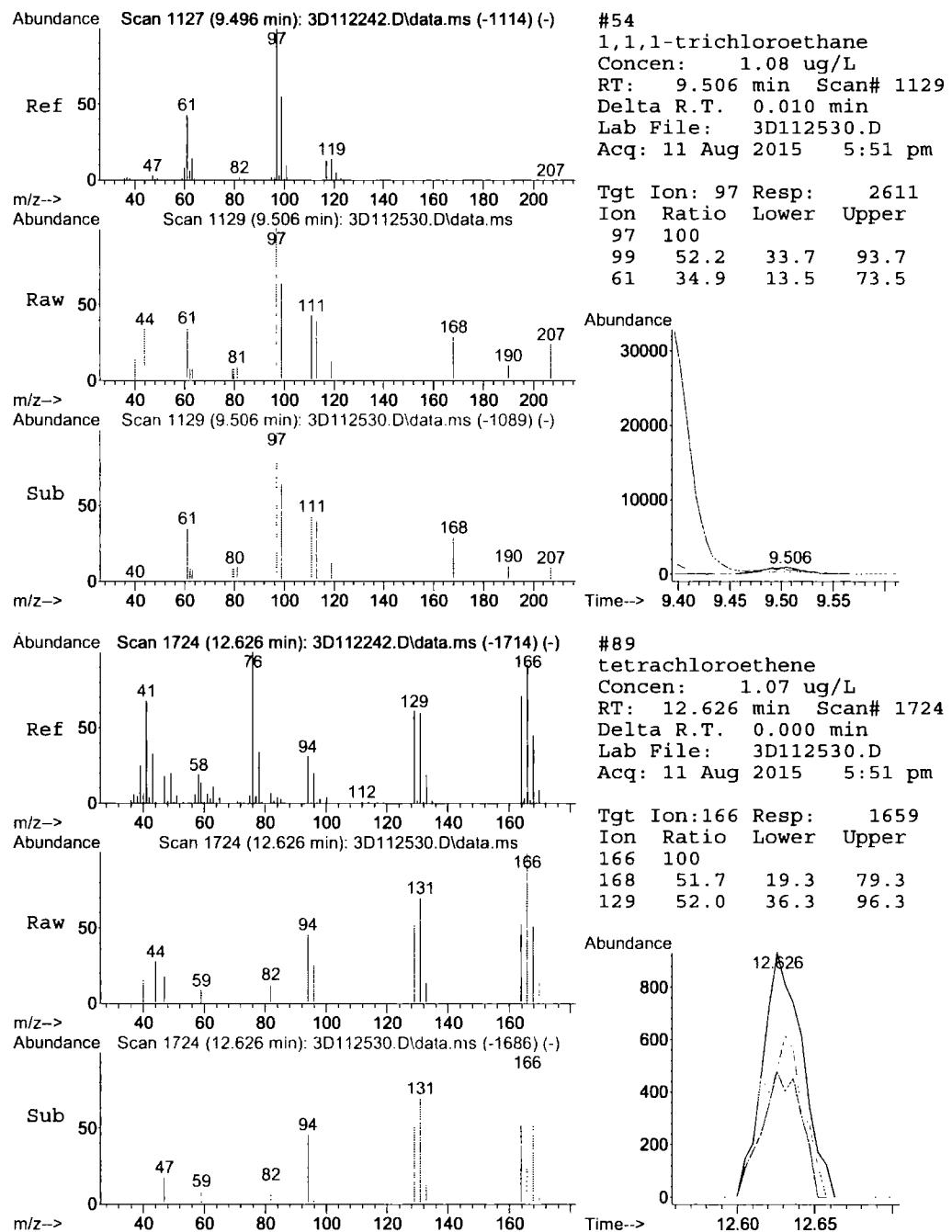
(#= qualifier out of range (m)= manual integration (+)= signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112530.D
 Acq On : 11 Aug 2015 5:51 pm
 Operator : ximenac
 Sample : jc1106-6
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 23 Sample Multiplier: 1

Quant Time: Aug 12 09:28:33 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112522.D
 Acq On : 11 Aug 2015 2:09 pm
 Operator : ximenac
 Sample : jc1106-7
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 15 Sample Multiplier: 1

Quant Time: Aug 11 16:11:14 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

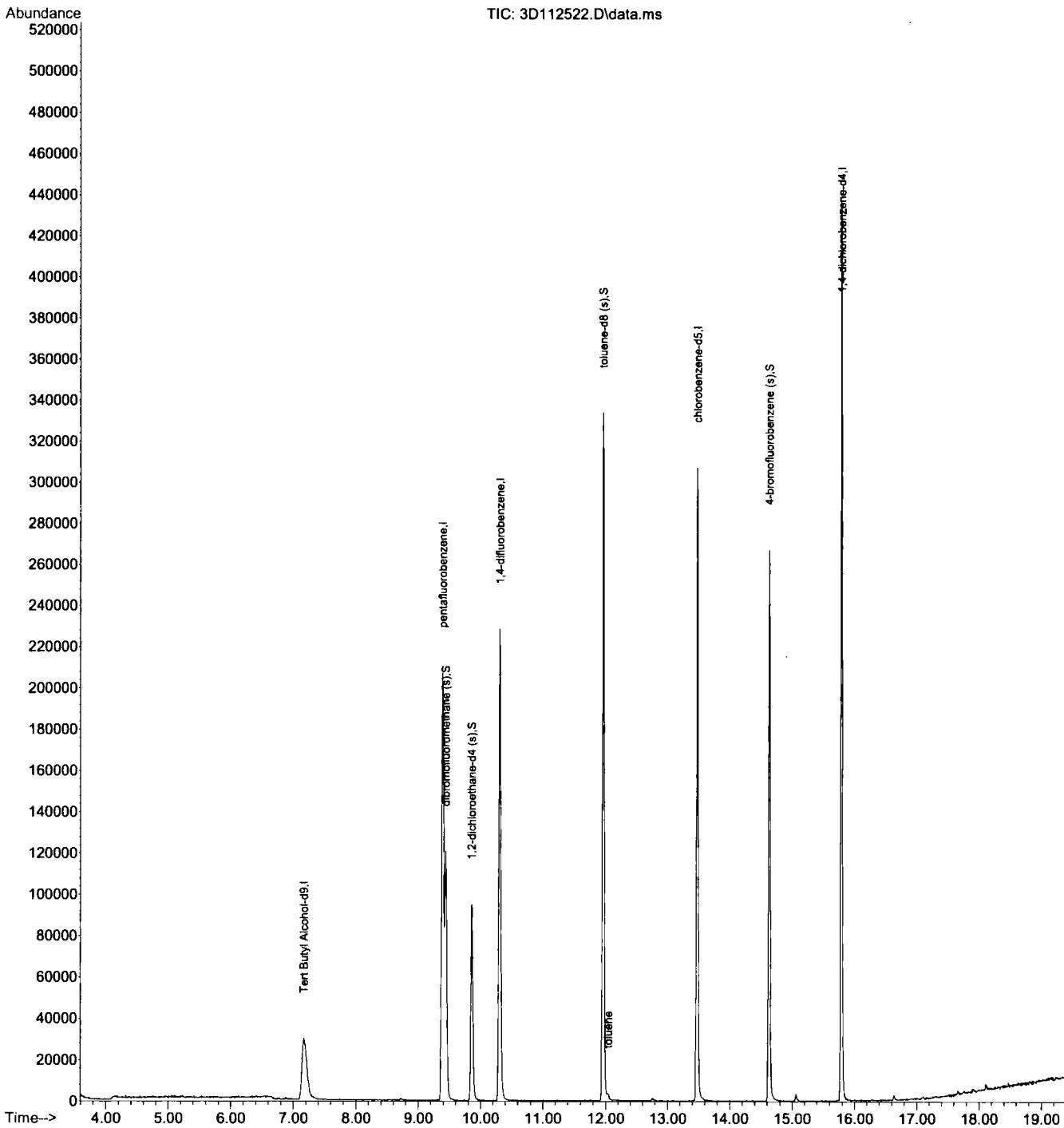
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.168	65	92150	500.00	ug/L	-0.01
4) pentafluorobenzene	9.391	168	192784	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	222789	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	184238	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	117577	50.00	ug/L	0.00
<hr/>						
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	82180	55.64	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	111.28%
51) 1,2-dichloroethane-d4 (s)	9.857	65	76411	55.84	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	111.68%
80) toluene-d8 (s)	11.960	98	245781	50.37	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	100.74%
106) 4-bromofluorobenzene (s)	14.629	95	86896	48.04	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.08%
<hr/>						
Target Compounds				Ovalue		
82) toluene	12.044	92	1283	0.45	ug/L	# 76

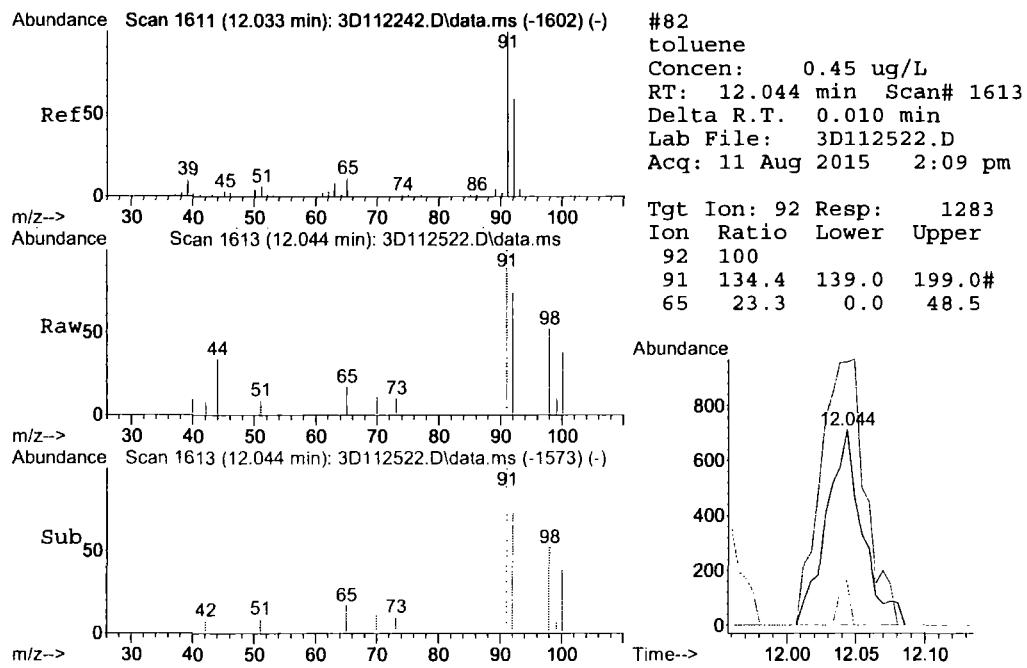
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112522.D
 Acq On : 11 Aug 2015 2:09 pm
 Operator : ximenac
 Sample : jc1106-7
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 15 Sample Multiplier: 1

Quant Time: Aug 11 16:11:14 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



7.17
7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\

Data File : 3D112535.D

Acq On : 11 Aug 2015 8:07 pm

Operator : ximenac

Sample : jc1106-8

Misc : MS89468,V3D4822,5,,,1

ALS Vial : 28 Sample Multiplier: 1

Quant Time: Aug 12 09:31:03 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

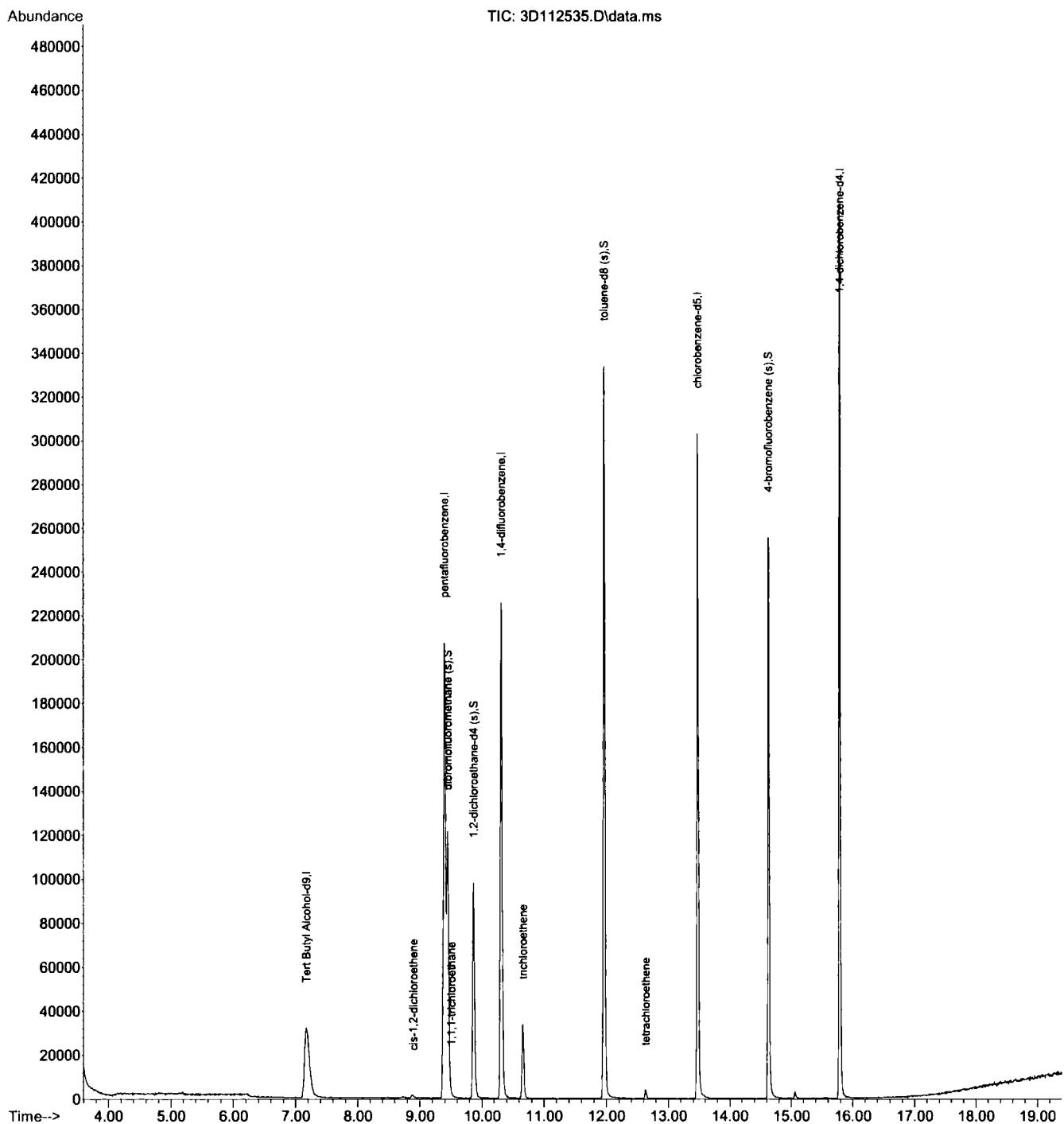
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	99250	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	189230	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	220375	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	181092	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	111525	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	82180	56.68	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	=	113.36%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	77091	57.40	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	=	114.80%	
80) toluene-d8 (s)	11.960	98	244029	50.56	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	=	101.12%	
106) 4-bromofluorobenzene (s)	14.629	95	82877	48.30	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	=	96.60%	
Target Compounds						
43) cis-1,2-dichloroethene	8.898	96	1150	0.70	ug/L	# 51
54) 1,1,1-trichloroethane	9.501	97	815	0.34	ug/L	86
69) trichloroethene	10.660	95	13296	10.62	ug/L	94
89) tetrachloroethene	12.631	166	1331	0.86	ug/L	86

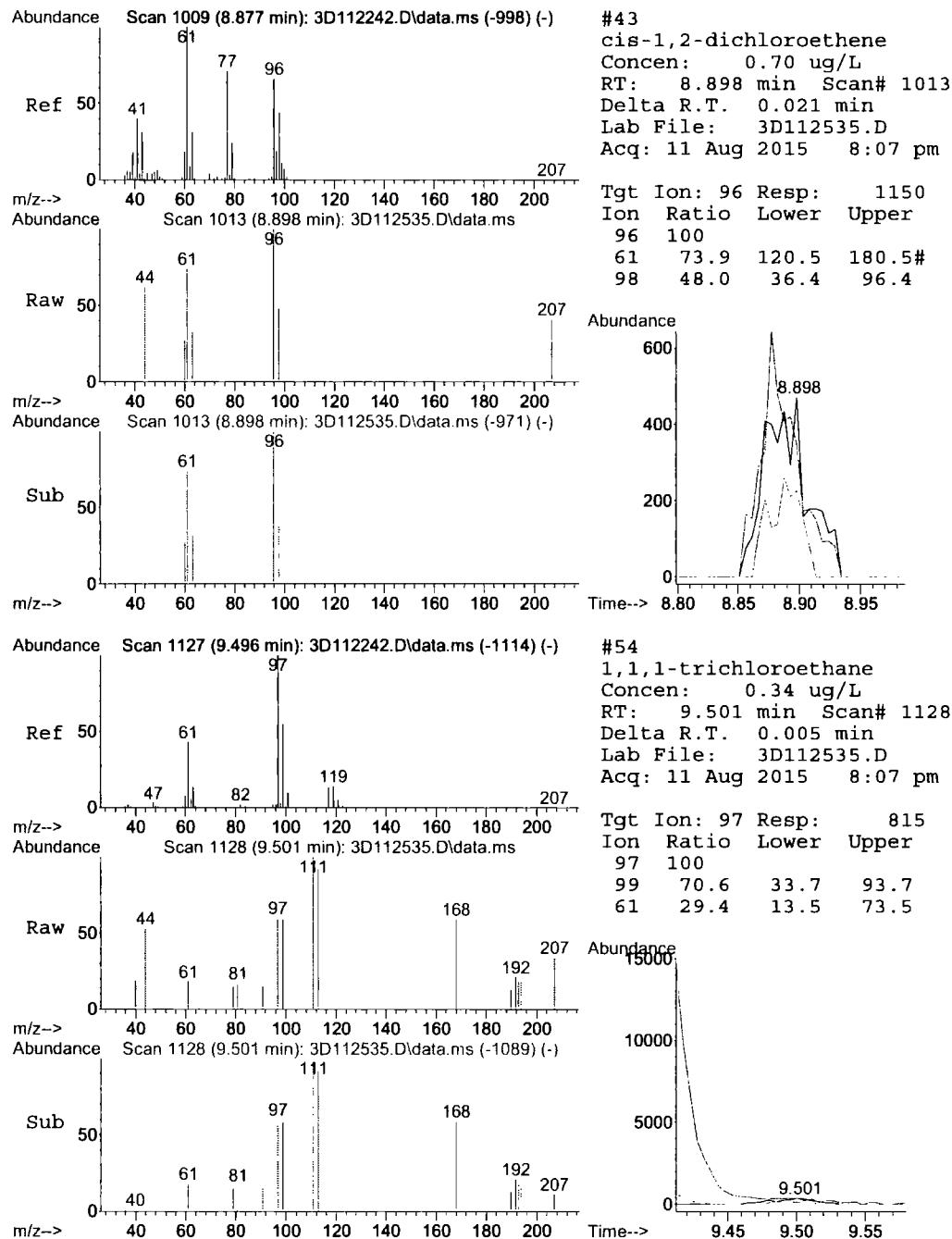
(#) = qualifier out of range (m) = manual integration (+) = signals summed

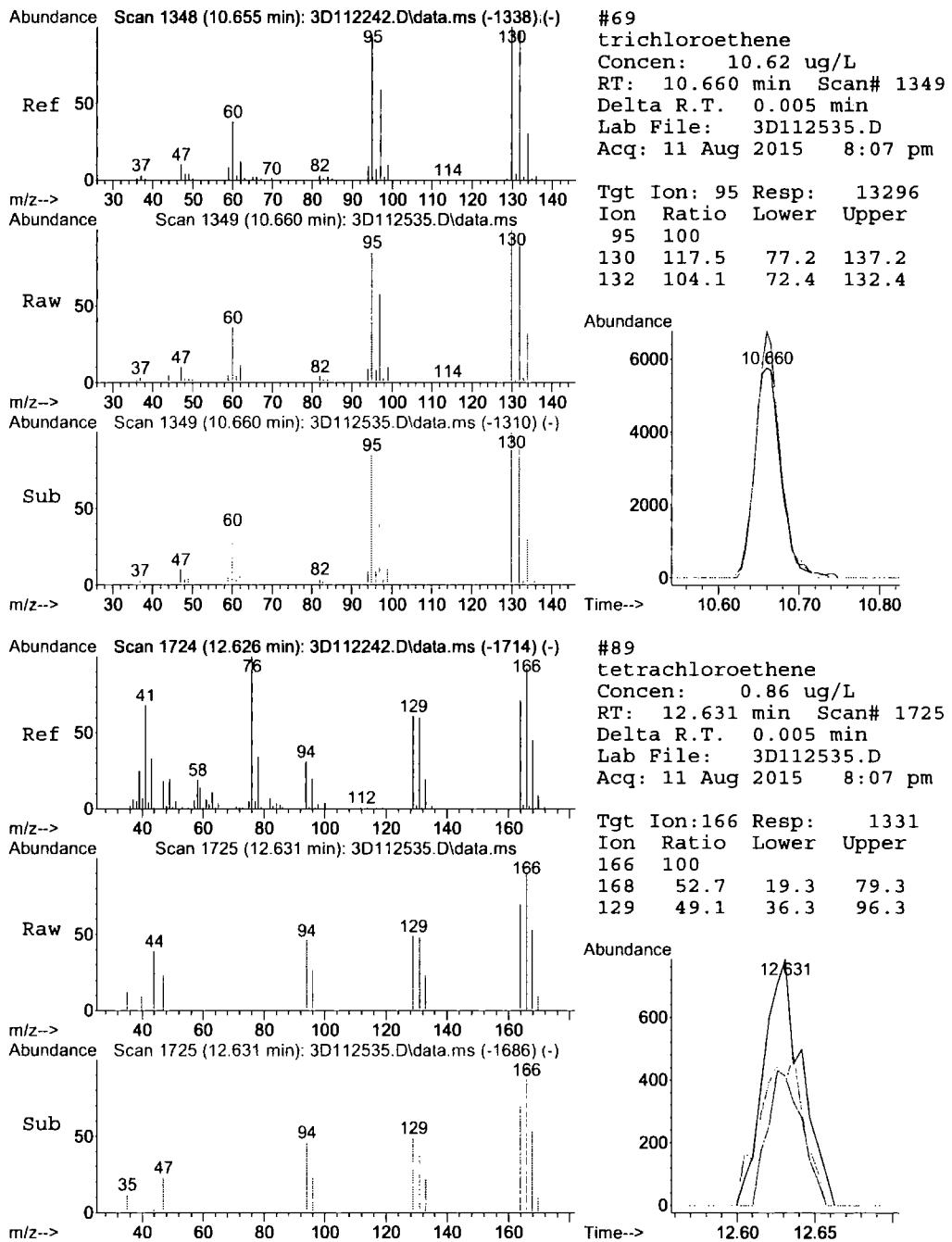
Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112535.D
 Acq On : 11 Aug 2015 8:07 pm
 Operator : ximenac
 Sample : jc1106-8
 Misc : MS89468, V3D4822,5,,1
 ALS Vial : 28 Sample Multiplier: 1

Quant Time: Aug 12 09:31:03 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration







Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\

Data File : 3D112572.D

Acq On : 12 Aug 2015 1:27 pm

Operator : ximenac

Sample : jc1106-9

Misc : MS89468,V3D4824,5,,,1

ALS Vial : 11 Sample Multiplier: 1

Quant Time: Aug 12 14:31:45 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

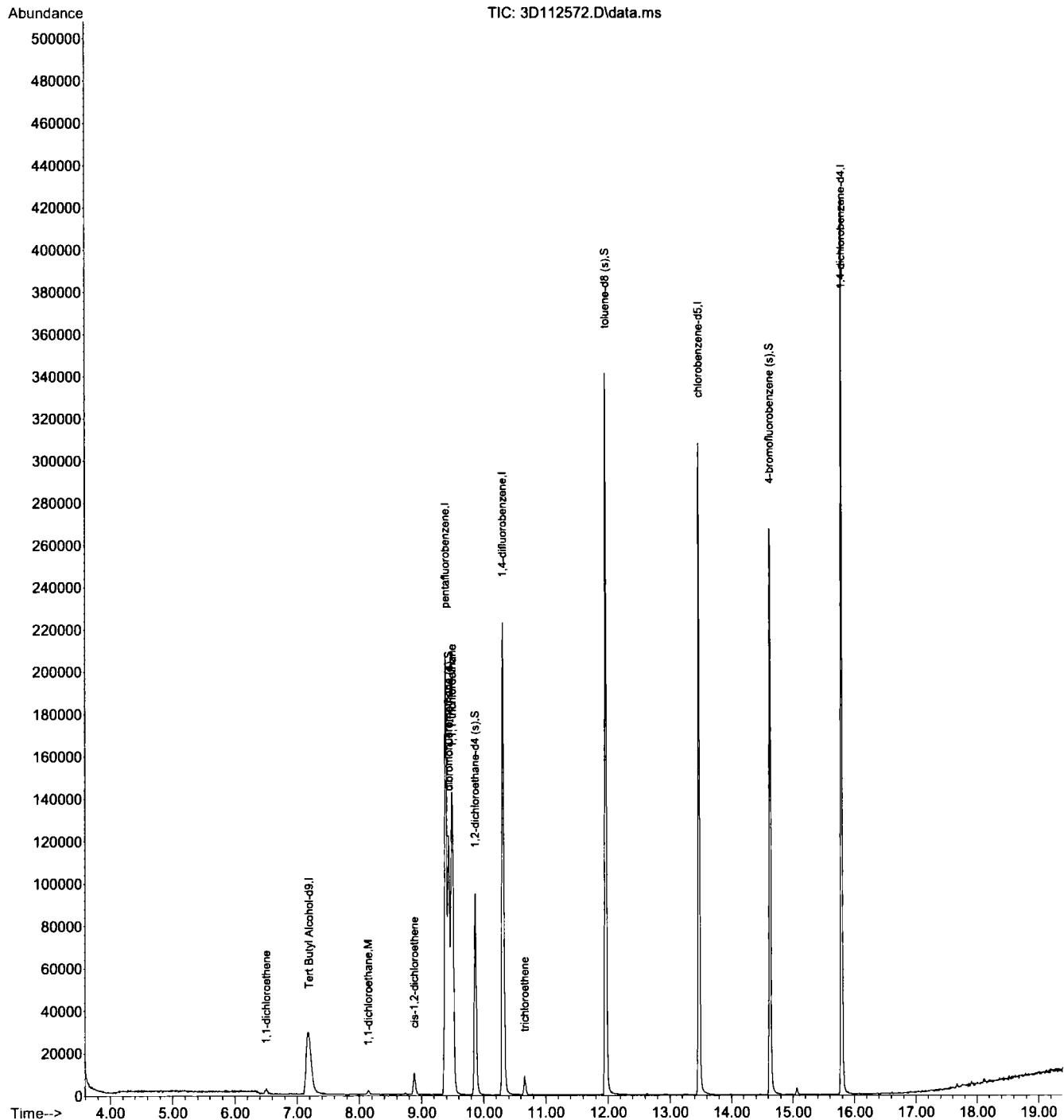
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.189	65	91436	500.00	ug/L	0.01
4) pentafluorobenzene	9.391	168	193182	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	221437	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	184042	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	114924	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	83193	56.21	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	112.42%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	76435	55.75	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	111.50%	
80) toluene-d8 (s)	11.960	98	247377	51.01	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	102.02%	
106) 4-bromofluorobenzene (s)	14.629	95	85271	48.22	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	96.44%	
Target Compounds						
				Qvalue		
22) 1,1-dichloroethene	6.502	61	2142	0.93	ug/L	93
37) 1,1-dichloroethane	8.138	63	2631	1.06	ug/L	93
43) cis-1,2-dichloroethene	8.882	96	5951	3.53	ug/L	# 79
54) 1,1,1-trichloroethane	9.496	97	132947	53.61	ug/L	96
69) trichloroethene	10.660	95	3265	2.60	ug/L	90

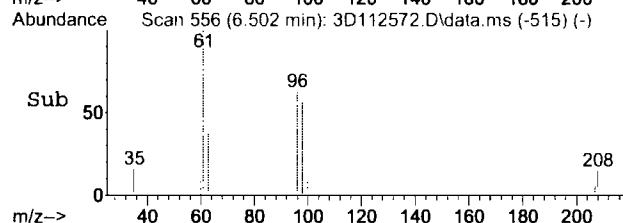
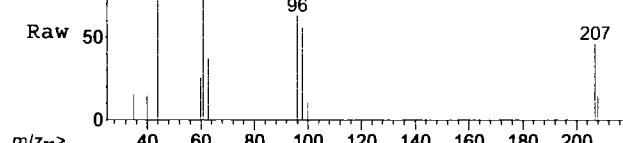
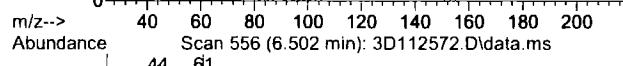
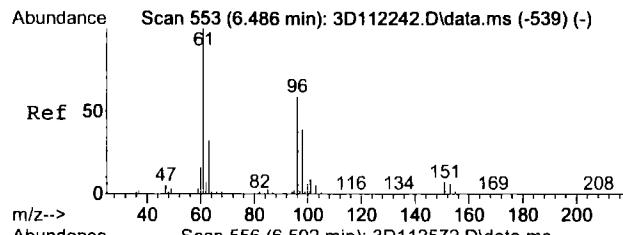
(#= qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
 Data File : 3D112572.D
 Acq On : 12 Aug 2015 1:27 pm
 Operator : ximenac
 Sample : jc1106-9
 Misc : MS89468,V3D4824,5,,,.1
 ALS Vial : 11 Sample Multiplier: 1

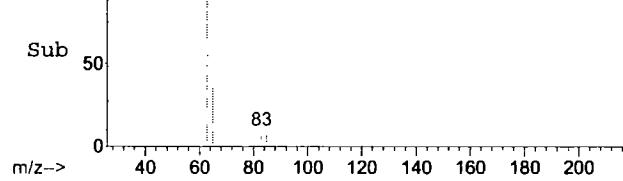
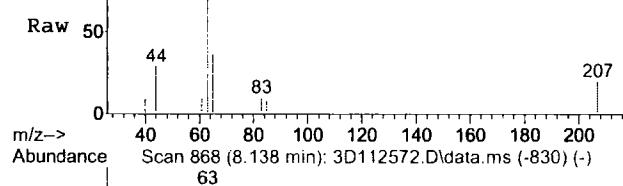
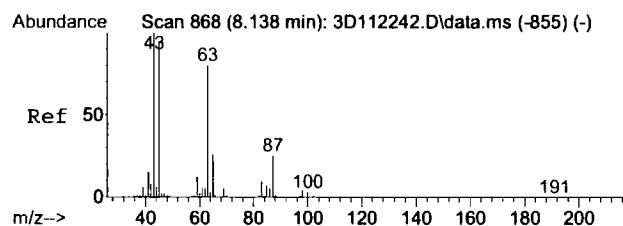
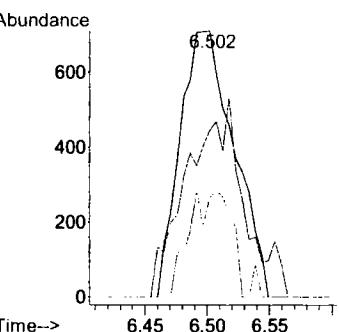
Quant Time: Aug 12 14:31:45 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration





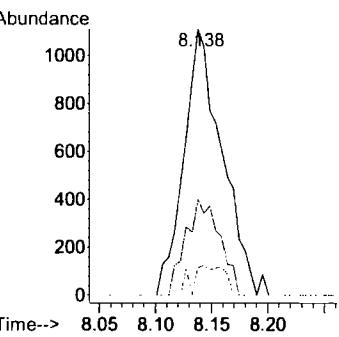
#22
1,1-dichloroethene
Concen: 0.93 ug/L
RT: 6.502 min Scan# 556
Delta R.T. 0.016 min
Lab File: 3D112572.D
Acq: 12 Aug 2015 1:27 pm

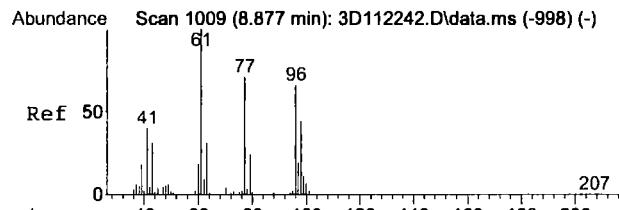
Tgt Ion: 61 Resp: 2142
Ion Ratio Lower Upper
61 100
96 62.6 28.7 88.7
63 37.4 1.8 61.8



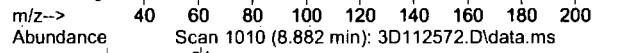
#37
1,1-dichloroethane
Concen: 1.06 ug/L
RT: 8.138 min Scan# 868
Delta R.T. 0.000 min
Lab File: 3D112572.D
Acq: 12 Aug 2015 1:27 pm

Tgt Ion: 63 Resp: 2631
Ion Ratio Lower Upper
63 100
65 36.3 2.4 62.4
83 10.4 0.0 42.8

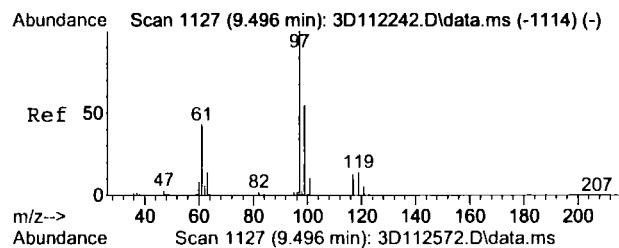
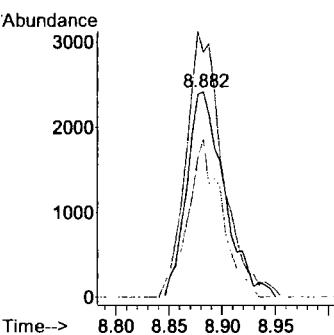
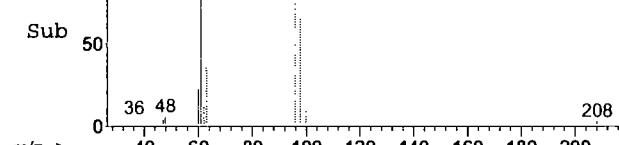
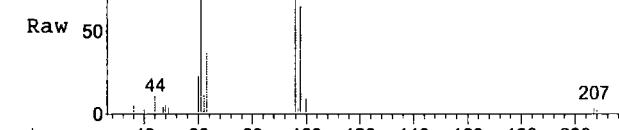




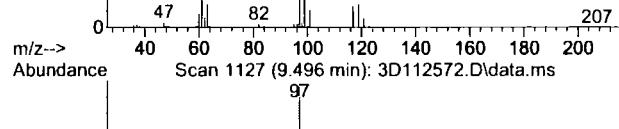
#43
cis-1,2-dichloroethene
Concen: 3.53 ug/L
RT: 8.882 min Scan# 1010
Delta R.T. 0.005 min
Lab File: 3D112572.D
Acq: 12 Aug 2015 1:27 pm



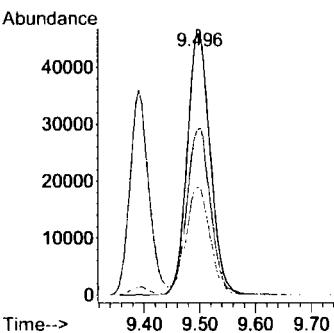
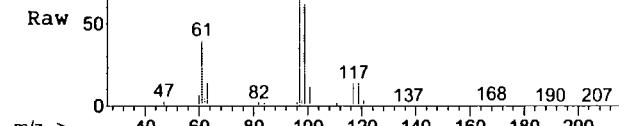
Tgt Ion: 96 Resp: 5951
Ion Ratio Lower Upper
96 100
61 119.1 120.5 180.5#
98 77.0 36.4 96.4

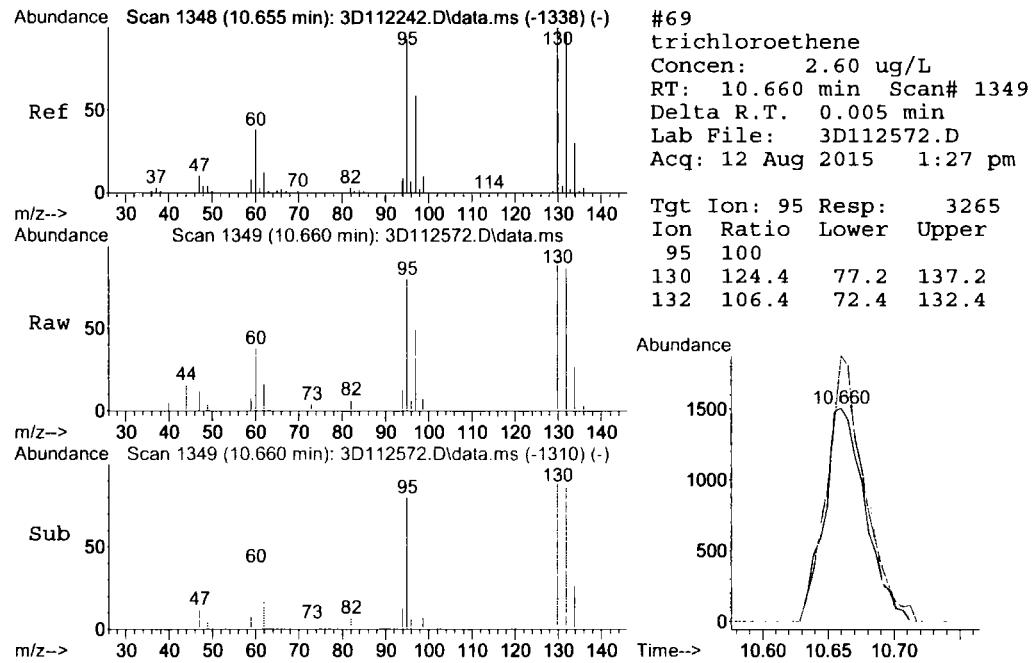


#54
1,1,1-trichloroethane
Concen: 53.61 ug/L
RT: 9.496 min Scan# 1127
Delta R.T. 0.000 min
Lab File: 3D112572.D
Acq: 12 Aug 2015 1:27 pm



Tgt Ion: 97 Resp: 132947
Ion Ratio Lower Upper
97 100
99 61.6 33.7 93.7
61 40.3 13.5 73.5





Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112525.D
 Acq On : 11 Aug 2015 3:30 pm
 Operator : ximenac
 Sample : jc1106-10
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 18 Sample Multiplier: 1

Quant Time: Aug 11 16:12:16 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.173	65	87229	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	179050	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.313	114	210954	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	172979	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	107549	50.00	ug/L	0.00

System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	77633	56.59	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	113.18%
51) 1,2-dichloroethane-d4 (s)	9.863	65	72314	56.90	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	113.80%
80) toluene-d8 (s)	11.960	98	229189	49.61	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	99.22%
106) 4-bromofluorobenzene (s)	14.629	95	79858	48.26	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.52%

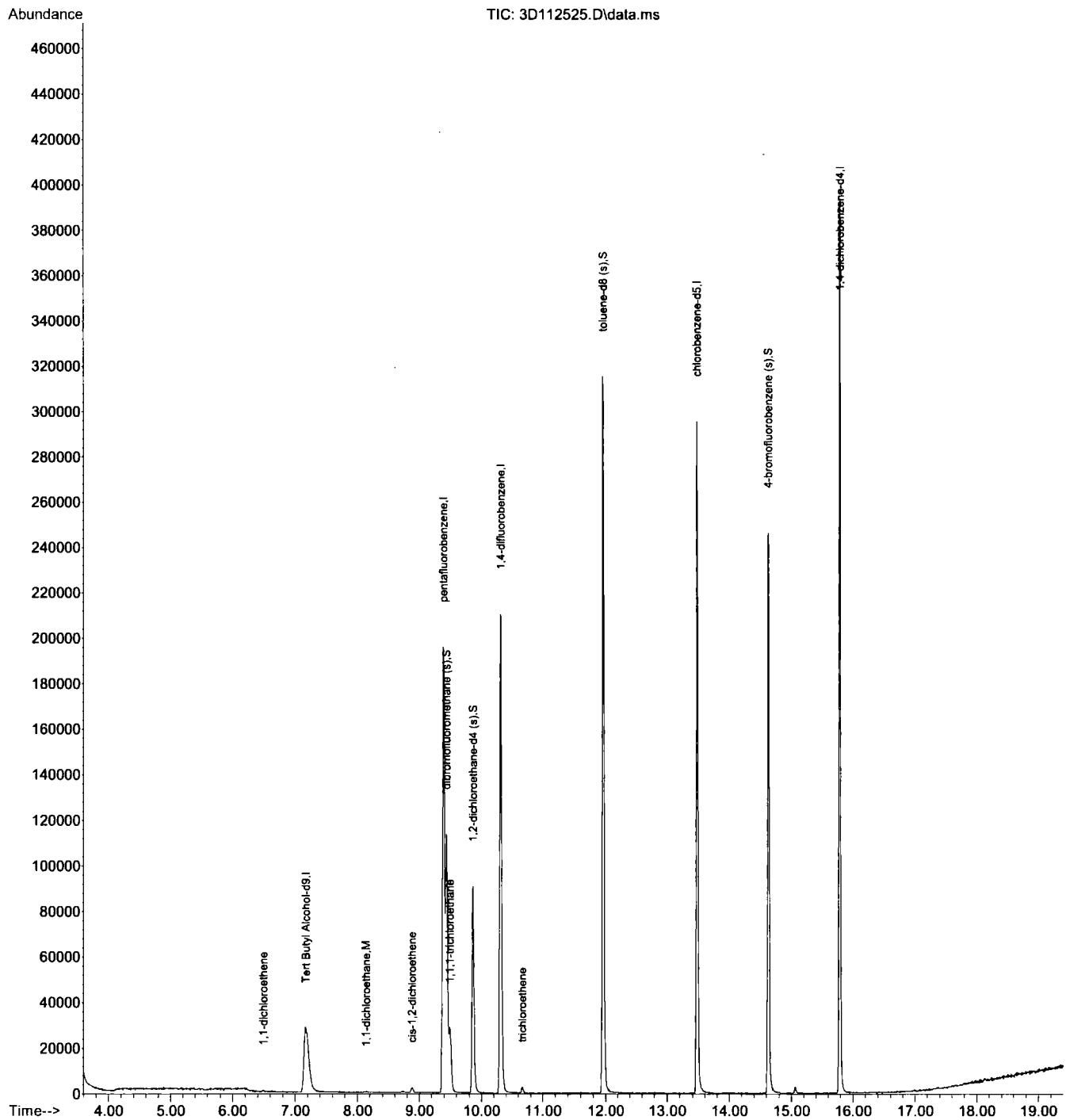
Target Compounds						Qvalue
22) 1,1-dichloroethene	6.497	61	434	0.20	ug/L	83
37) 1,1-dichloroethane	8.143	63	581	0.25	ug/L	# 49
43) cis-1,2-dichloroethene	8.877	96	1541	0.99	ug/L	# 69
54) 1,1,1-trichloroethane	9.496	97	22968	9.99	ug/L	94
69) trichloroethene	10.670	95	1004	0.84	ug/L	84

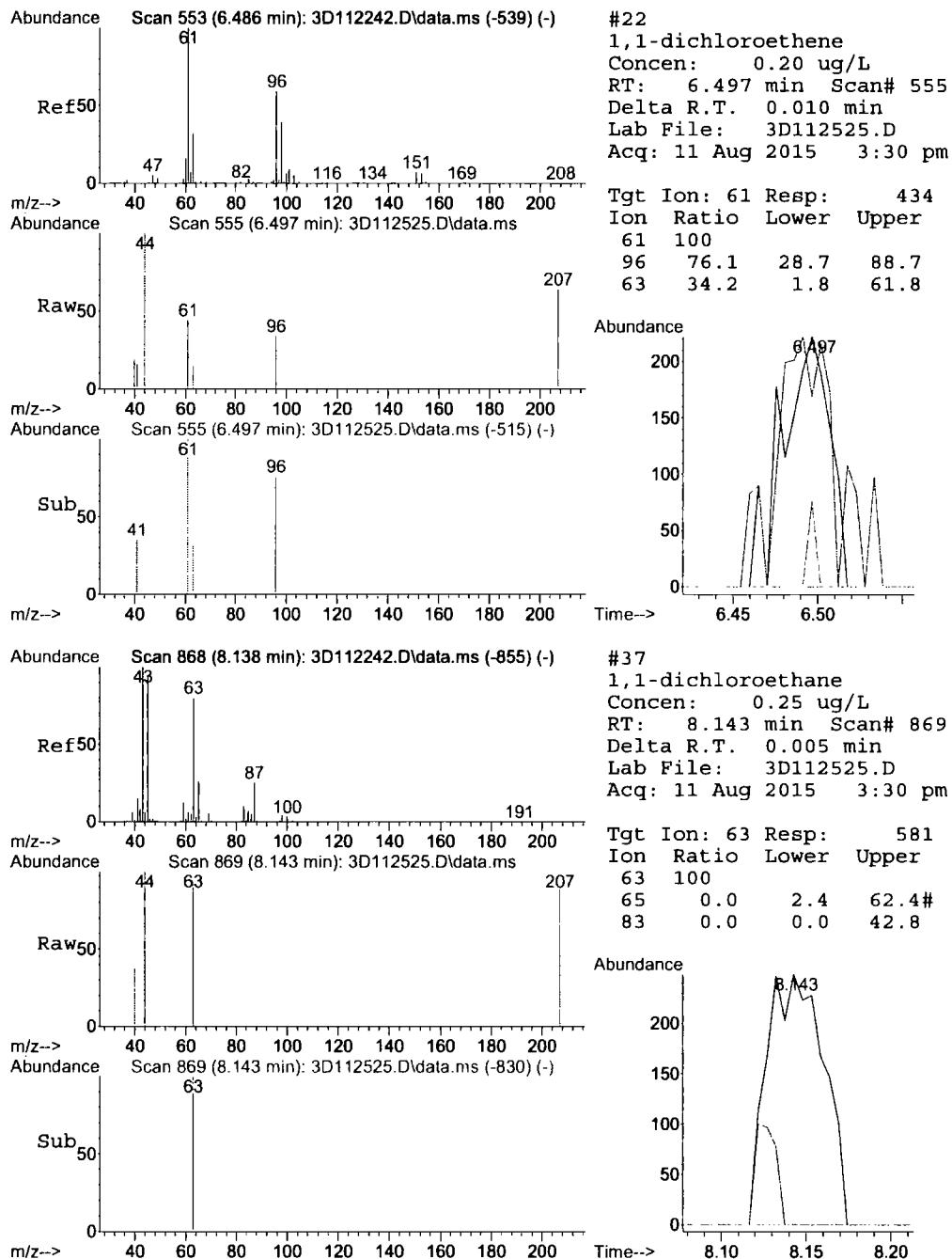
(#) = qualifier out of range (m) = manual integration (+) = signals summed

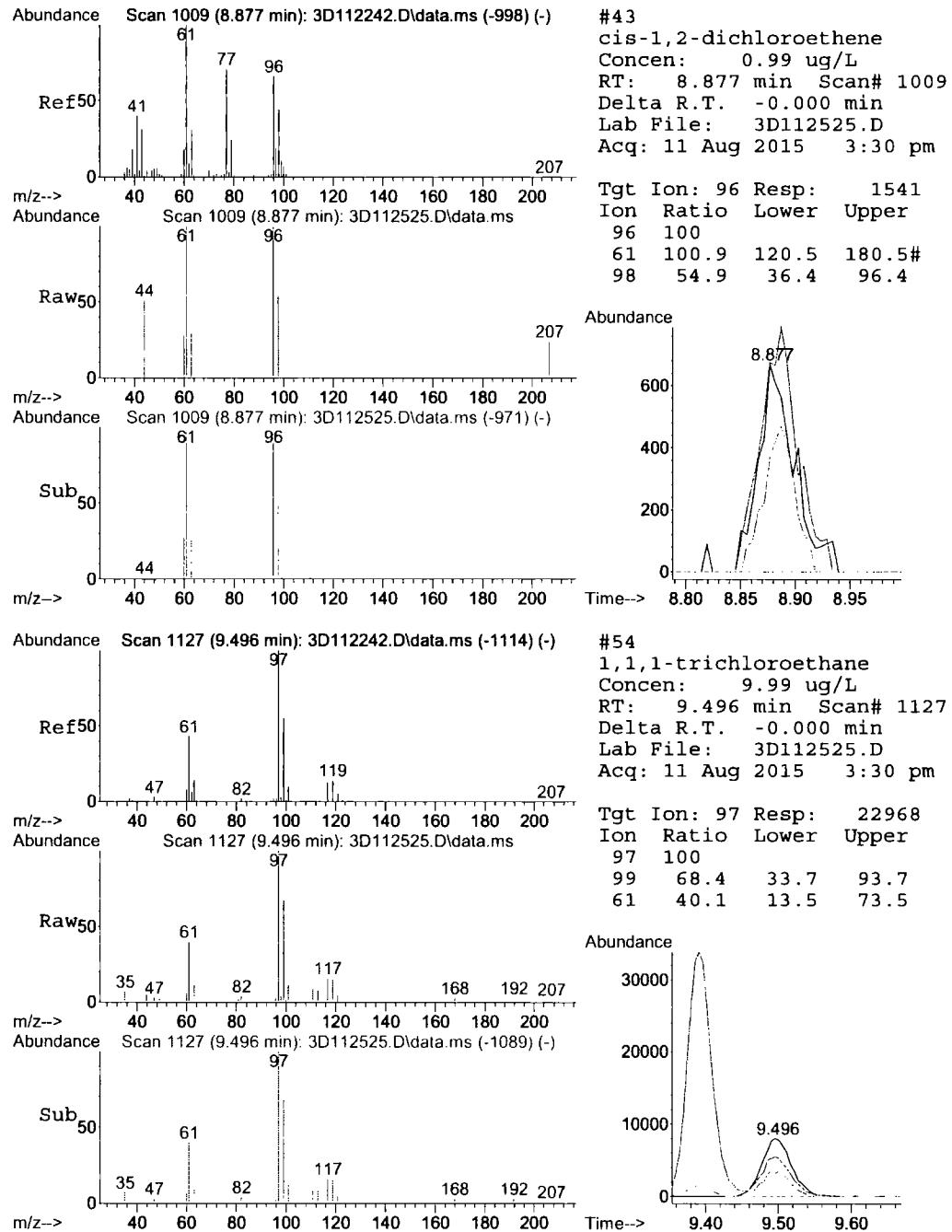
Quantitation Report (QT Reviewed)

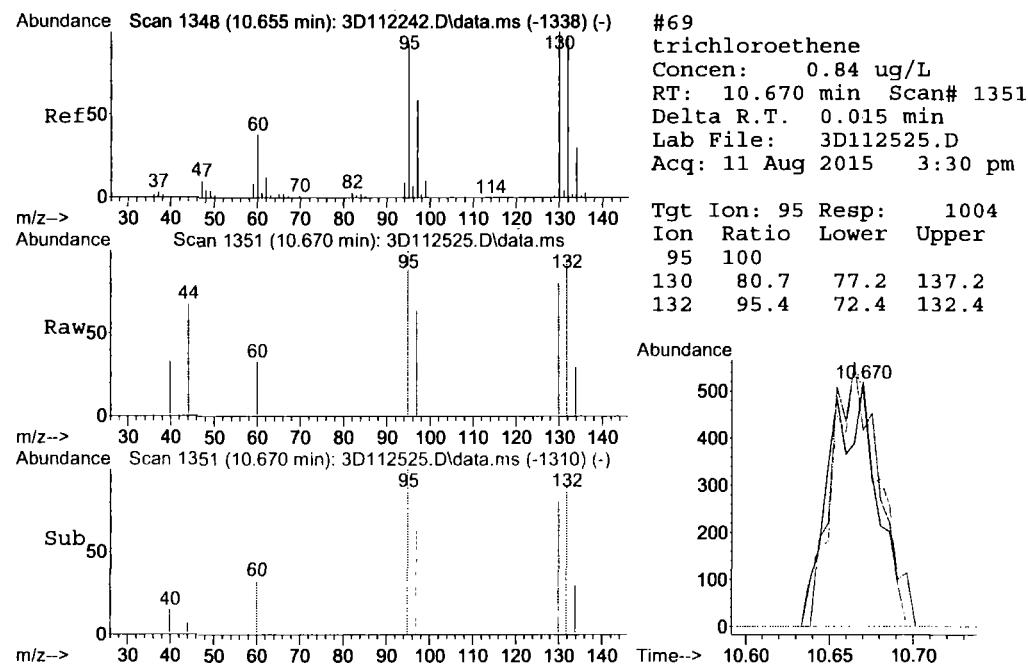
Data Path : C:\msdchem\1\DATA\
 Data File : 3D112525.D
 Acq On : 11 Aug 2015 3:30 pm
 Operator : ximenac
 Sample : jc1106-10
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 18 Sample Multiplier: 1

Quant Time: Aug 11 16:12:16 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration









Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112538.D
 Acq On : 11 Aug 2015 9:28 pm
 Operator : ximenac
 Sample : jc1106-11
 Misc : MS89468,V3D4822,5,,,,1
 ALS Vial : 31 Sample Multiplier: 1

Quant Time: Aug 12 09:31:55 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
<hr/>						
Internal Standards						
1) Tert Butyl Alcohol-d9	7.168	65	84171	500.00	ug/L	-0.01
4) pentafluorobenzene	9.391	168	181857	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.313	114	212162	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	174691	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.792	152	107285	50.00	ug/L	0.00
<hr/>						
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.438	113	79322	56.93	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	113.86%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	72709	56.33	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	112.66%	
80) toluene-d8 (s)	11.960	98	234906	50.55	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	101.10%	
106) 4-bromofluorobenzene (s)	14.629	95	80040	48.49	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	96.98%	

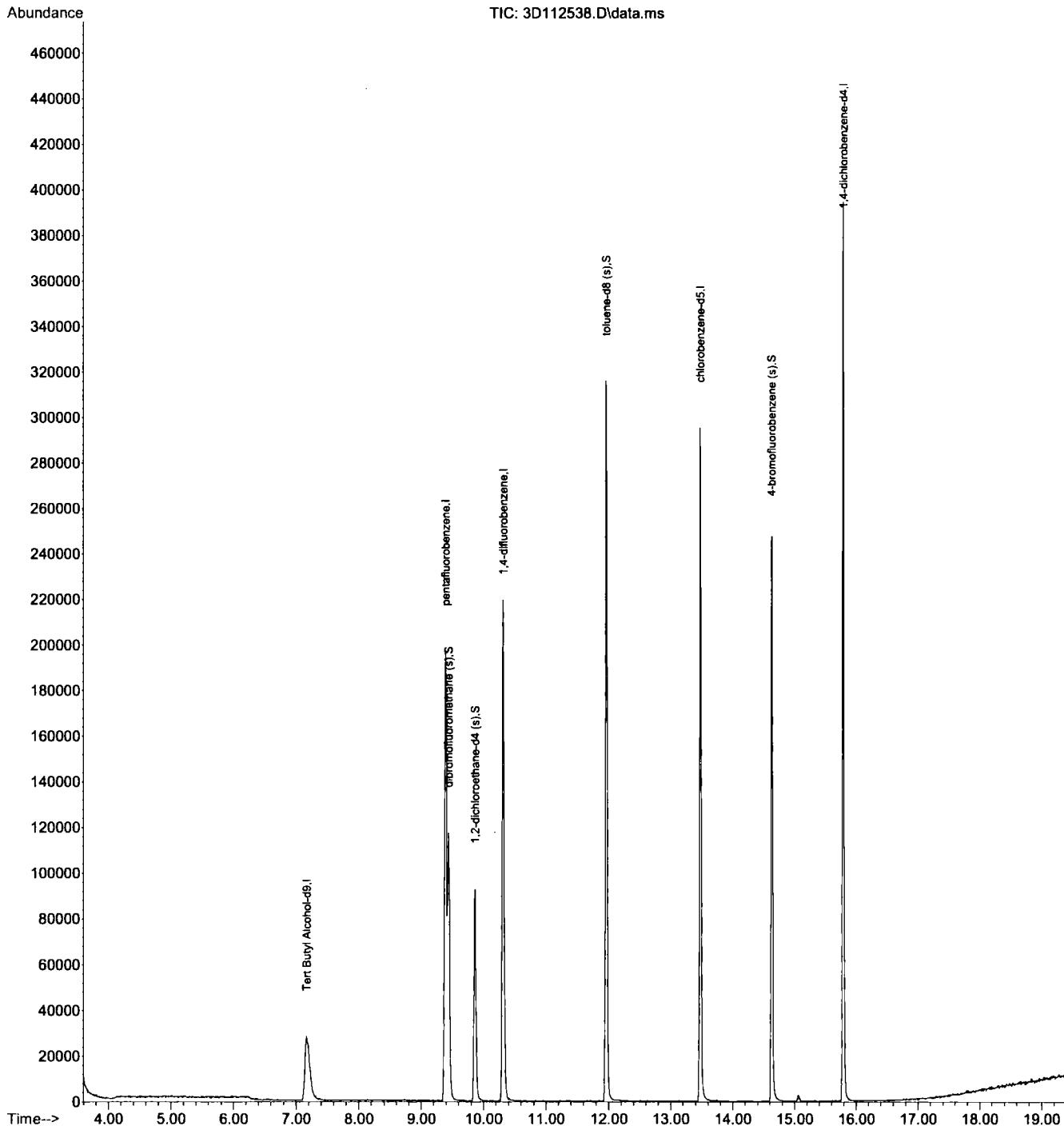
Target Compounds	Qvalue
<hr/>	

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
Data File : 3D112538.D
Acq On : 11 Aug 2015 9:28 pm
Operator : ximenac
Sample : jc1106-11
Misc : MS89468,V3D4822,5,,,1
ALS Vial : 31 Sample Multiplier: 1

Quant Time: Aug 12 09:31:55 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\

Data File : 3D112537.D

Acq On : 11 Aug 2015 9:01 pm

Operator : ximenac

Sample : jc1106-12

Misc : MS89468,V3D4822,5,,,1

ALS Vial : 30 Sample Multiplier: 1

Quant Time: Aug 12 09:31:42 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

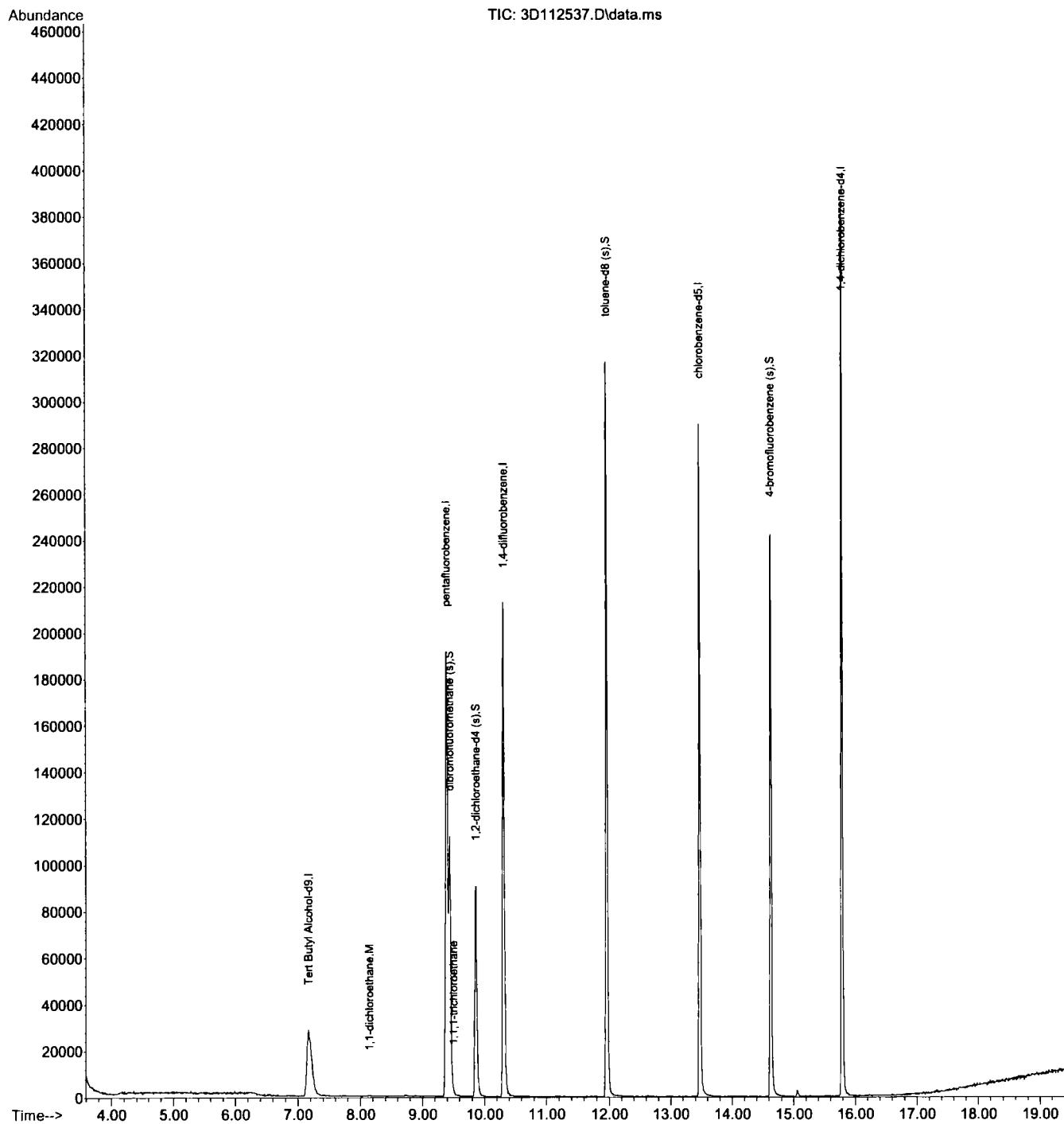
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.168	65	84872	500.00	ug/L	-0.01
4) pentafluorobenzene	9.391	168	176303	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	208560	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	169036	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	106176	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	76587	56.70	ug/L	0.00
Spiked Amount 50.000 Range	76 - 120		Recovery	=	113.40%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	71854	57.42	ug/L	0.00
Spiked Amount 50.000 Range	73 - 122		Recovery	=	114.84%	
80) toluene-d8 (s)	11.960	98	226231	49.53	ug/L	0.00
Spiked Amount 50.000 Range	84 - 119		Recovery	=	99.06%	
106) 4-bromofluorobenzene (s)	14.629	95	78825	48.25	ug/L	0.00
Spiked Amount 50.000 Range	78 - 117		Recovery	=	96.50%	
Target Compounds						
				Value		
37) 1,1-dichloroethane	8.143	63	626	0.28 ug/L	82	
54) 1,1,1-trichloroethane	9.501	97	789	0.35 ug/L	#	68

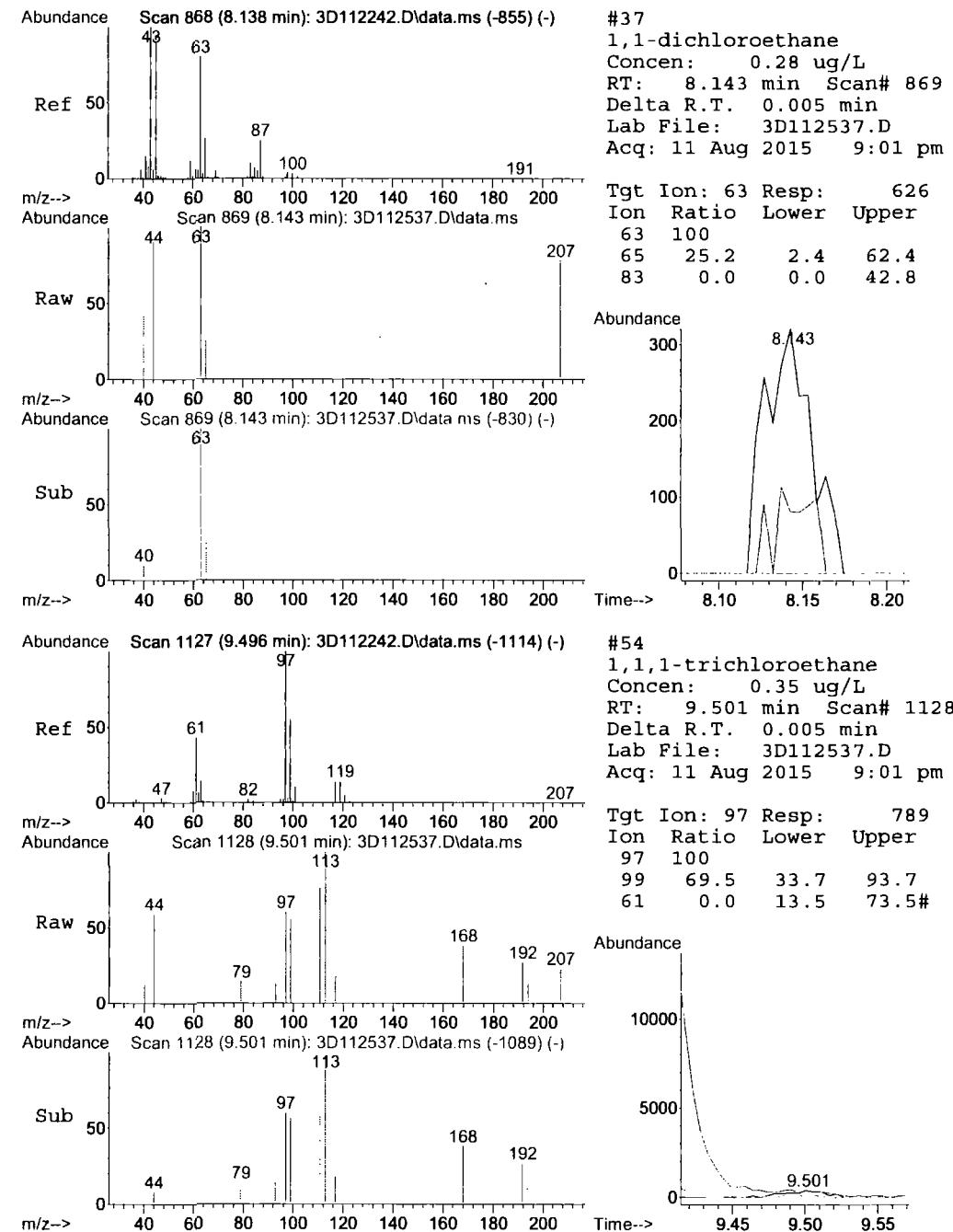
(#= qualifier out of range (m)= manual integration (+)= signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112537.D
 Acq On : 11 Aug 2015 9:01 pm
 Operator : ximenac
 Sample : jc1106-12
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 30 Sample Multiplier: 1

Quant Time: Aug 12 09:31:42 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112524.D
 Acq On : 11 Aug 2015 3:03 pm
 Operator : ximenac
 Sample : jc1106-13
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Aug 11 16:11:51 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.173	65	91266	500.00	ug/L	0.00
4) pentafluorobenzene	9.396	168	186566	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	219744	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	177389	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	110292	50.00	ug/L	0.00

System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	79189	55.40	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	110.80%
51) 1,2-dichloroethane-d4 (s)	9.863	65	73979	55.87	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	111.74%
80) toluene-d8 (s)	11.960	98	236581	49.16	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	98.32%
106) 4-bromofluorobenzene (s)	14.629	95	82401	48.56	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.12%

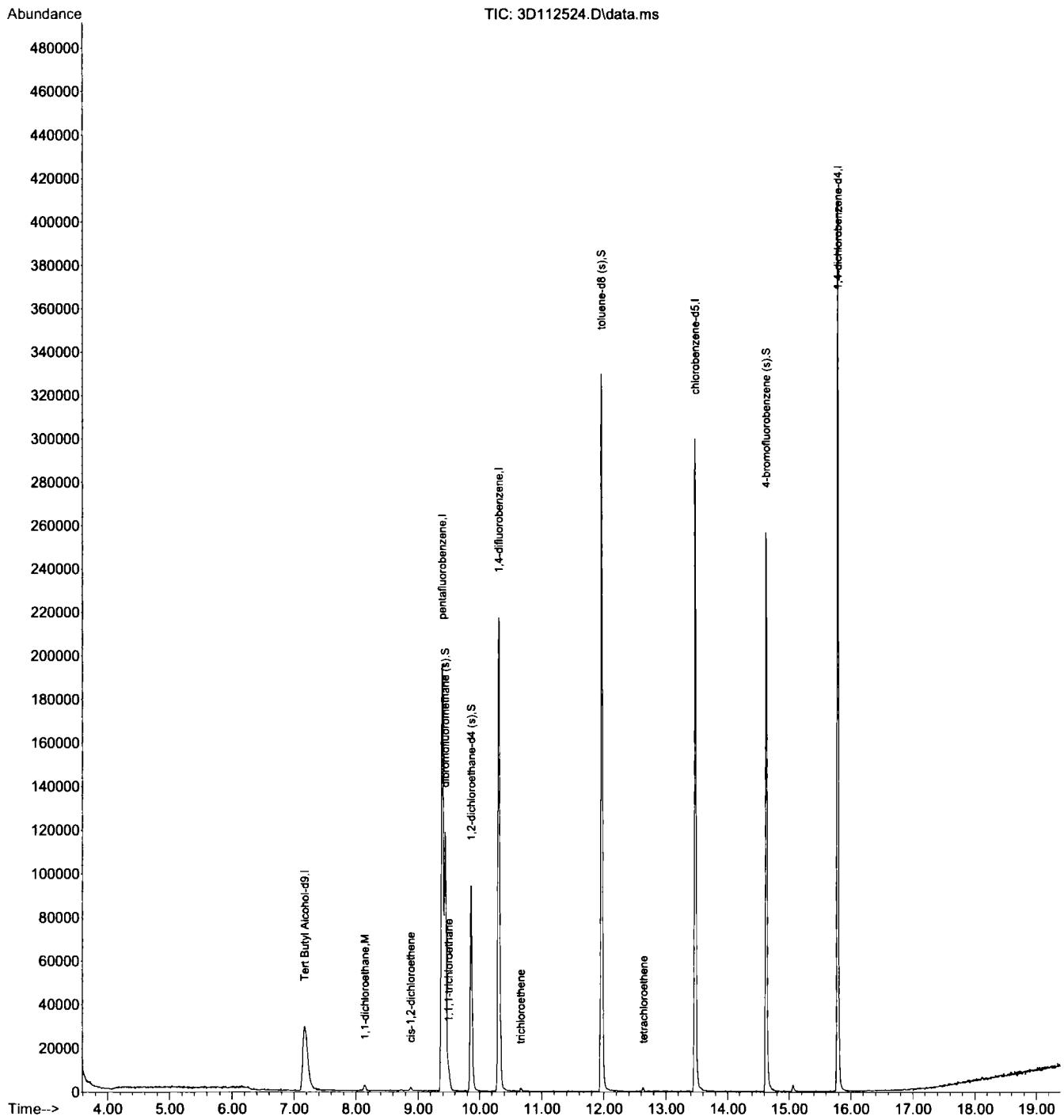
Target Compounds					Qvalue
37) 1,1-dichloroethane	8.143	63	3804	1.58	ug/L 87
43) cis-1,2-dichloroethene	8.882	96	921	0.57	ug/L 82
54) 1,1,1-trichloroethane	9.501	97	7121	2.97	ug/L 98
69) trichloroethene	10.665	95	498	0.40	ug/L # 58
89) tetrachloroethene	12.642	166	655	0.43	ug/L 71

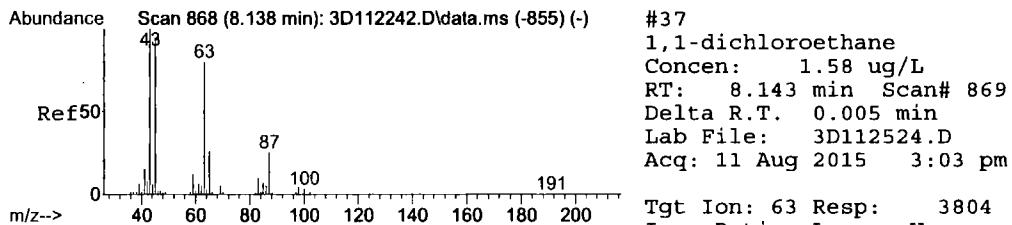
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

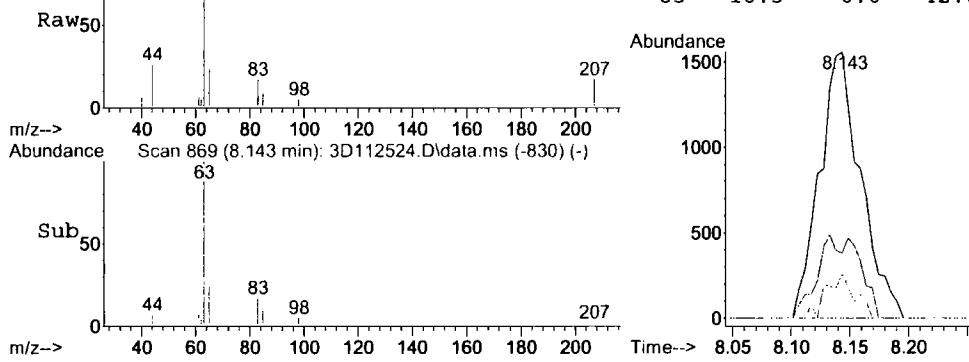
Data Path : C:\msdchem\1\DATA\
 Data File : 3D112524.D
 Acq On : 11 Aug 2015 3:03 pm
 Operator : ximenac
 Sample : jc1106-13
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Aug 11 16:11:51 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



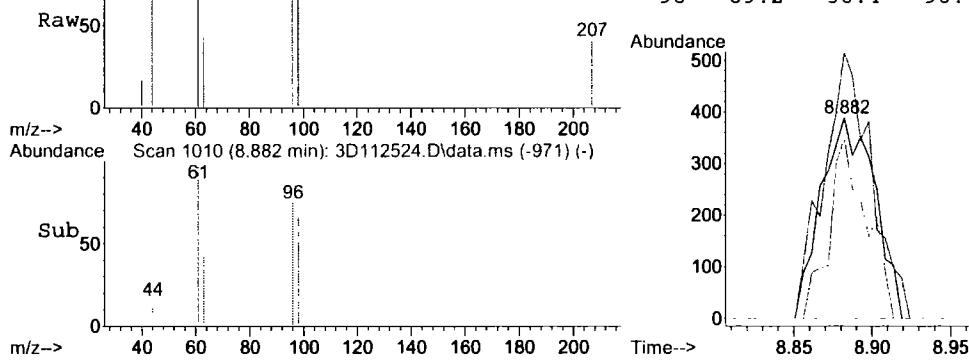


Tgt Ion: 63 Resp: 3804
Ion Ratio Lower Upper
63 100
65 24.5 2.4 62.4
83 16.5 0.0 42.8



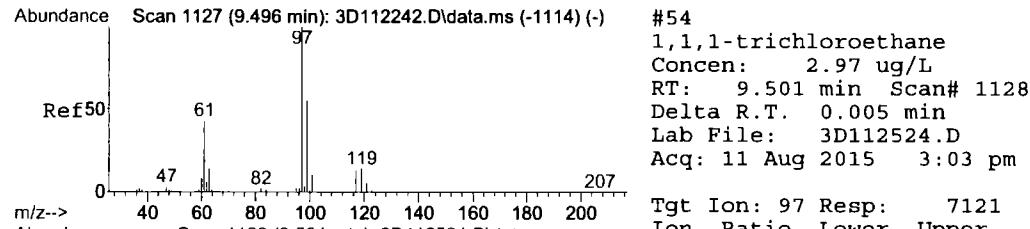
#43
cis-1,2-dichloroethene
Concen: 0.57 ug/L
RT: 8.882 min Scan# 1010
Delta R.T. 0.005 min
Lab File: 3D112524.D
Acq: 11 Aug 2015 3:03 pm

Tgt Ion: 96 Resp: 921
Ion Ratio Lower Upper
96 100
61 132.6 120.5 180.5
98 89.2 36.4 96.4

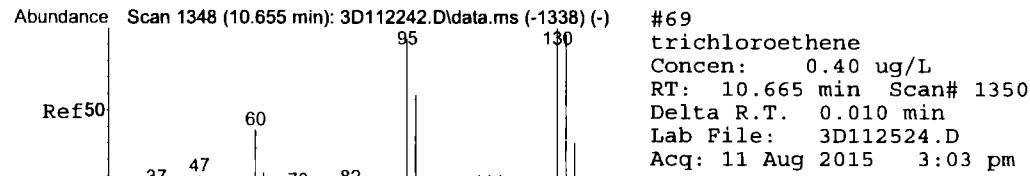
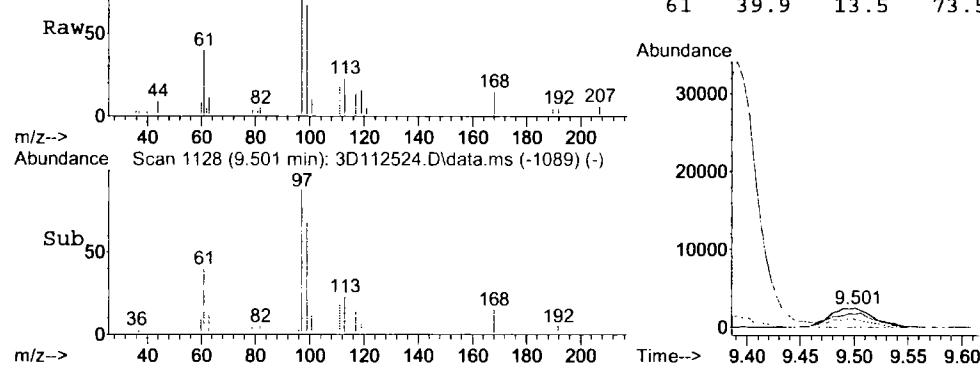


7.1.13

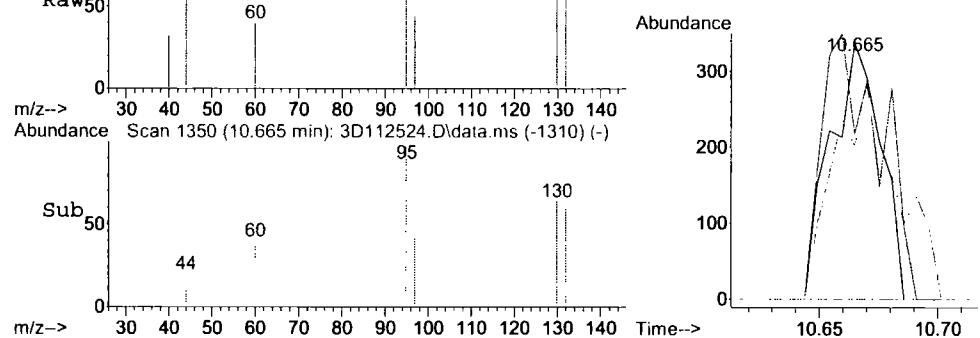
7

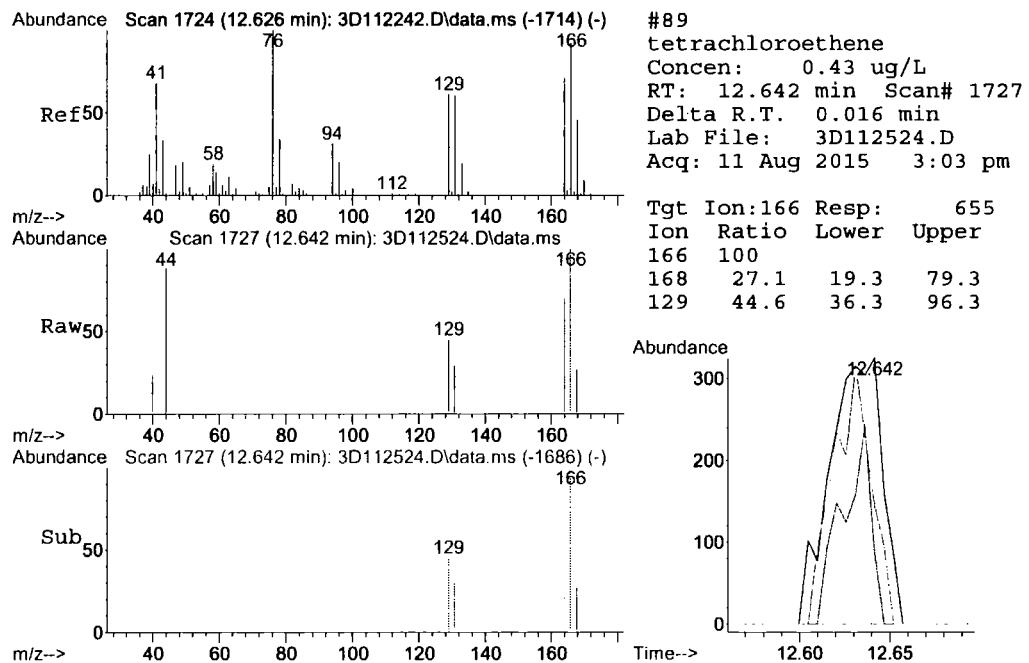


Tgt	Ion:	97	Resp:	7121
Ion	Ratio	Lower	Upper	
97	100			
99	63.8	33.7	93.7	
61	39.9	13.5	73.5	



Tgt	Ion:	95	Resp:	498
Ion	Ratio	Lower	Upper	
95	100			
130	64.5	77.2	137.2#	
132	59.8	72.4	132.4#	





7.1.13

7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\

Data File : 3D112536.D

Acq On : 11 Aug 2015 8:34 pm

Operator : ximenac

Sample : jc1106-14

Misc : MS89468,V3D4822,5,,,1

ALS Vial : 29 Sample Multiplier: 1

Quant Time: Aug 12 09:31:26 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

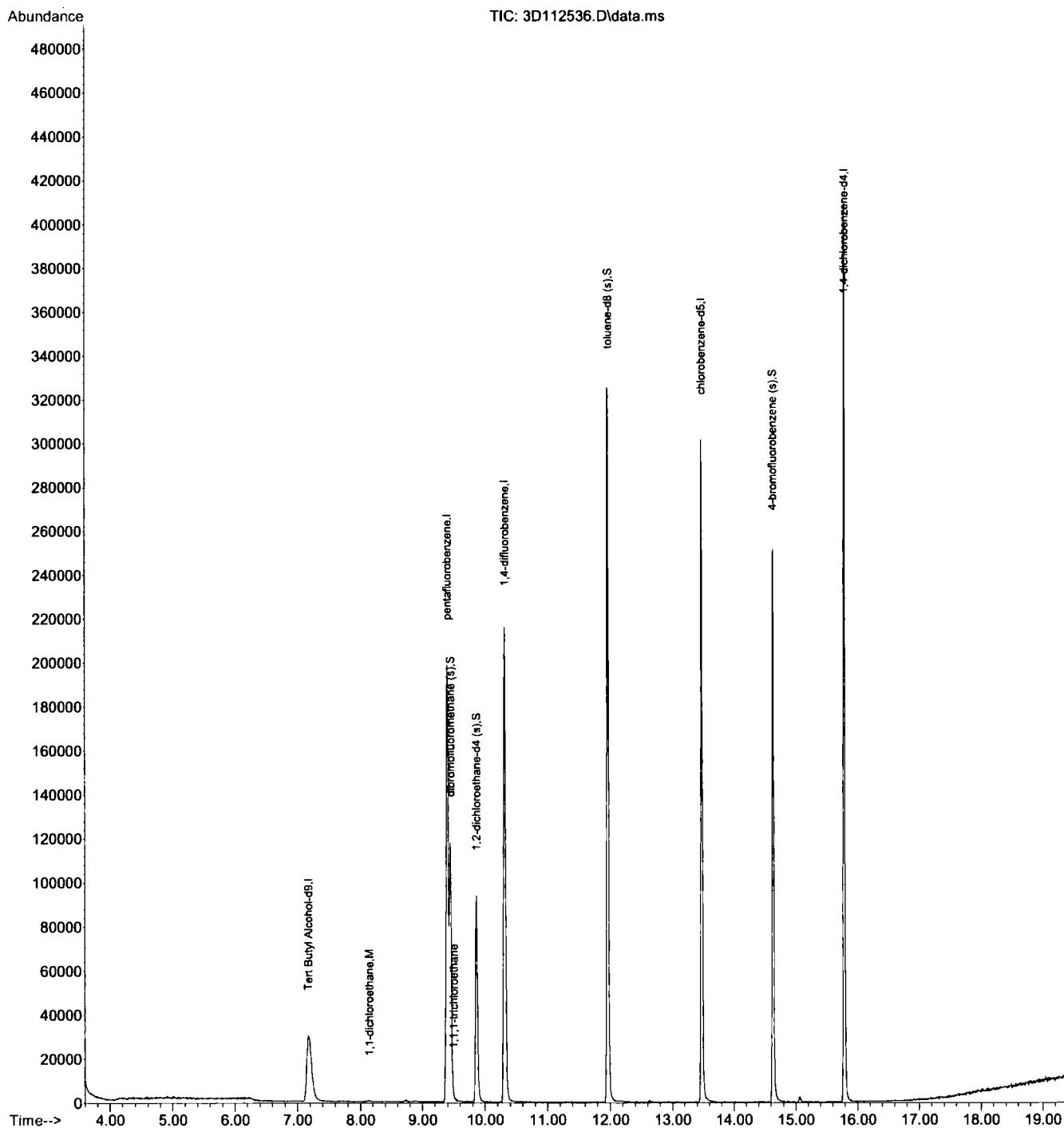
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	91107	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	183583	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	216133	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	177301	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	109805	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	80342	57.12	ug/L	0.00
Spiked Amount 50.000 Range	76 - 120		Recovery	=	114.24%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	75935	58.28	ug/L	0.00
Spiked Amount 50.000 Range	73 - 122		Recovery	=	116.56%	
80) toluene-d8 (s)	11.965	98	239928	50.69	ug/L	0.00
Spiked Amount 50.000 Range	84 - 119		Recovery	=	101.38%	
106) 4-bromofluorobenzene (s)	14.629	95	81846	48.45	ug/L	0.00
Spiked Amount 50.000 Range	78 - 117		Recovery	=	96.90%	
Target Compounds						
37) 1,1-dichloroethane	8.143	63	612	0.26	ug/L	# 49
54) 1,1,1-trichloroethane	9.496	97	845	0.36	ug/L	# 45

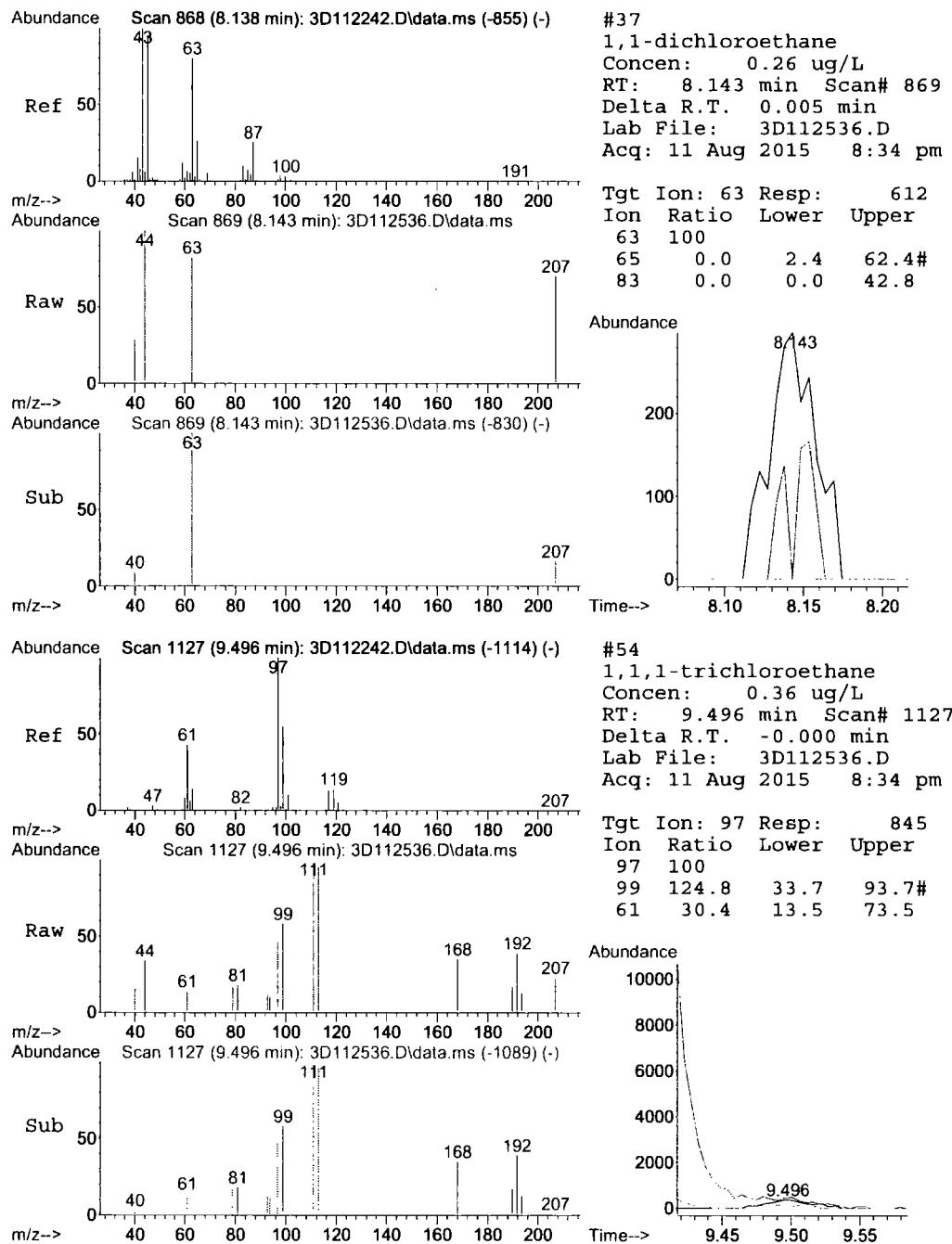
(#= qualifier out of range (m)= manual integration (+)= signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112536.D
 Acq On : 11 Aug 2015 8:34 pm
 Operator : ximenac
 Sample : jc1106-14
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 29 Sample Multiplier: 1

Quant Time: Aug 12 09:31:26 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112533.D
 Acq On : 11 Aug 2015 7:13 pm
 Operator : ximenac
 Sample : jc1106-15
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 26 Sample Multiplier: 1

Quant Time: Aug 12 09:29:30 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

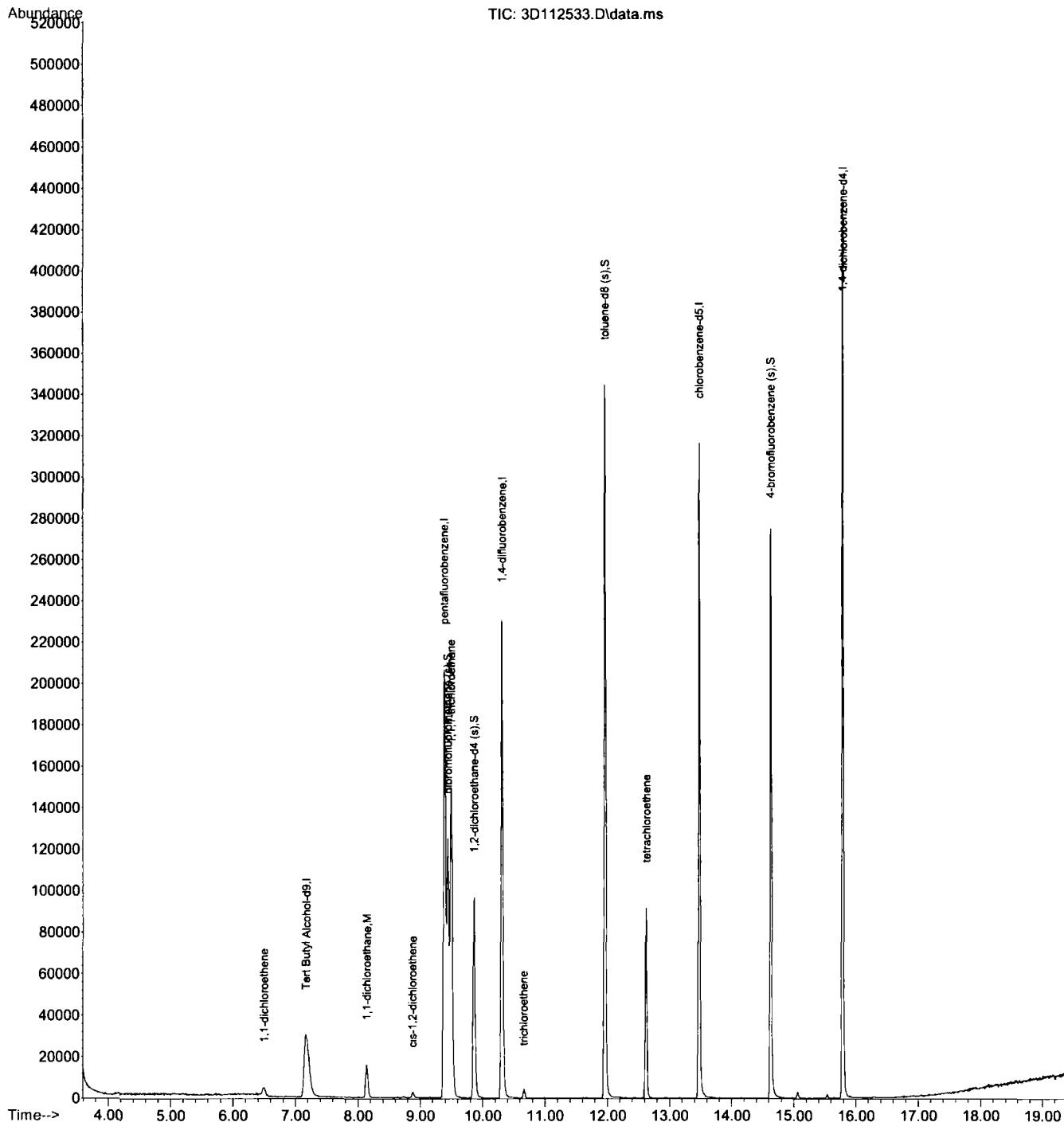
Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	94051	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	193794	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.313	114	228111	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	188105	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	116304	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	83552	56.27	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	112.54%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	77136	56.08	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	112.16%	
80) toluene-d8 (s)	11.960	98	250629	50.17	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	100.34%	
106) 4-bromofluorobenzene (s)	14.629	95	86098	48.11	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	96.22%	
Target Compounds						
				Qvalue		
22) 1,1-dichloroethene	6.497	61	4214	1.83	ug/L	98
37) 1,1-dichloroethane	8.143	63	21368	8.55	ug/L	98
43) cis-1,2-dichloroethene	8.877	96	1855	1.10	ug/L	80
54) 1,1,1-trichloroethane	9.496	97	135997	54.67	ug/L	99
69) trichloroethene	10.660	95	1664	1.28	ug/L	82
89) tetrachloroethene	12.626	166	30763	19.21	ug/L	95

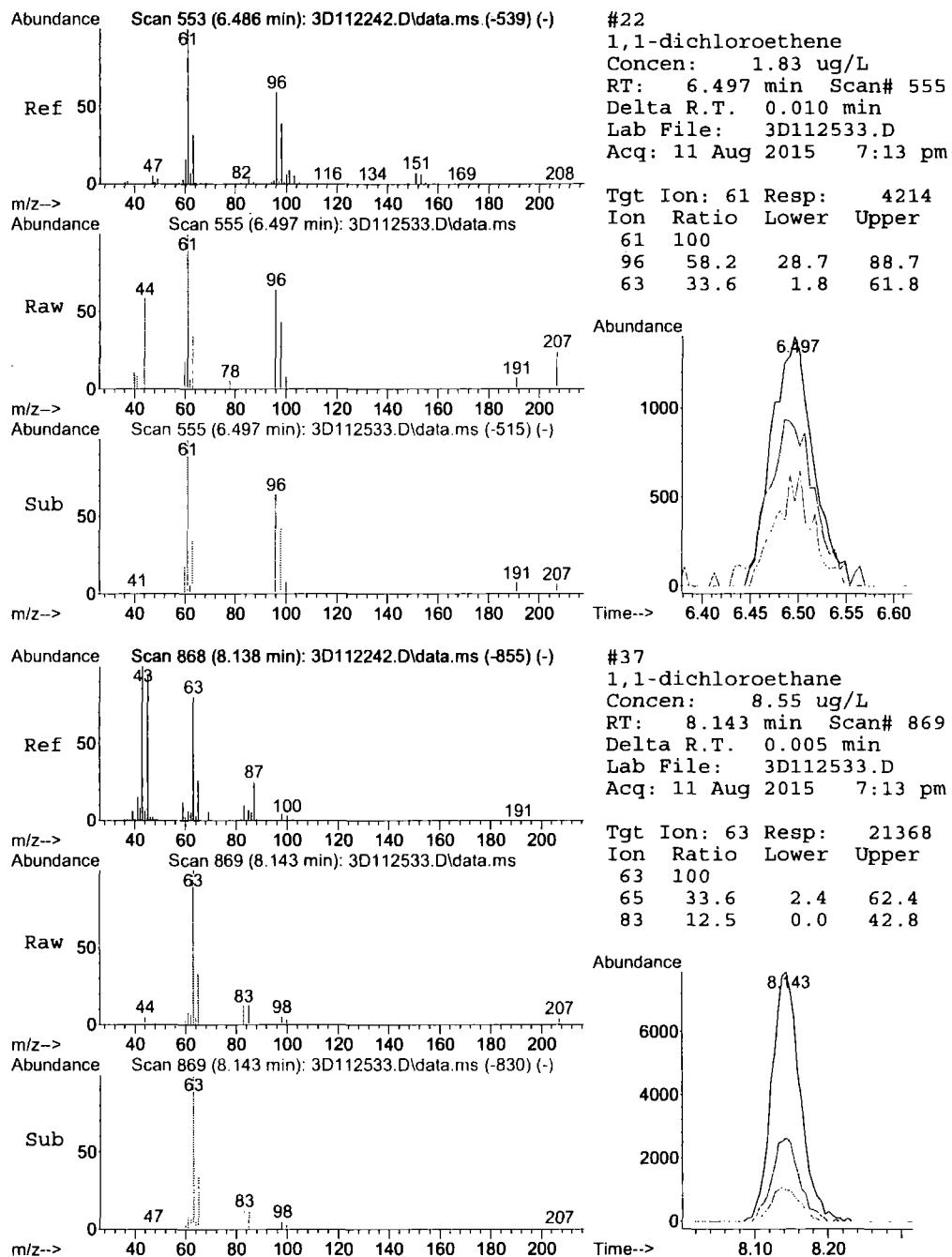
(#) = qualifier out of range (m) = manual integration (+) = signals summed

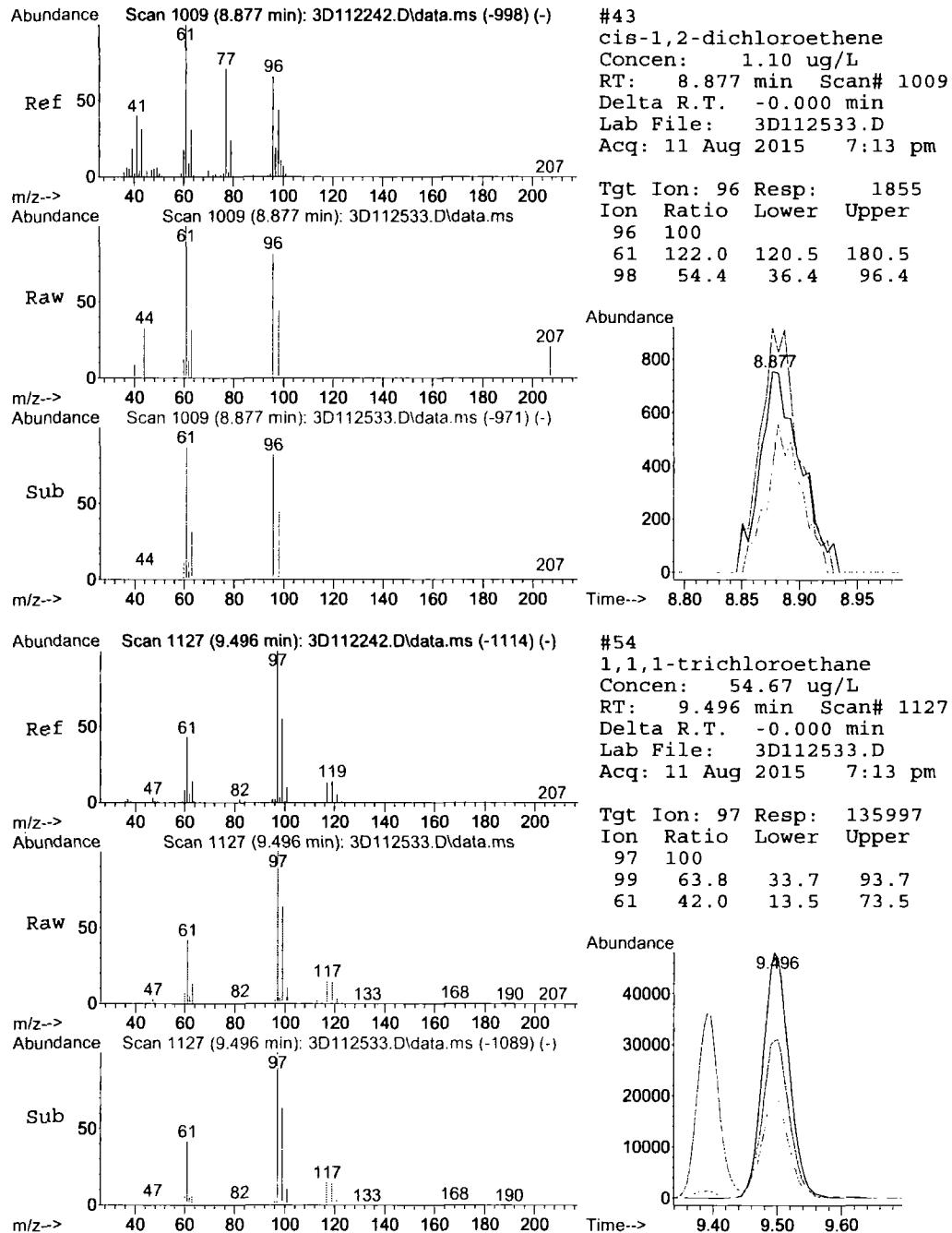
Quantitation Report (QT Reviewed)

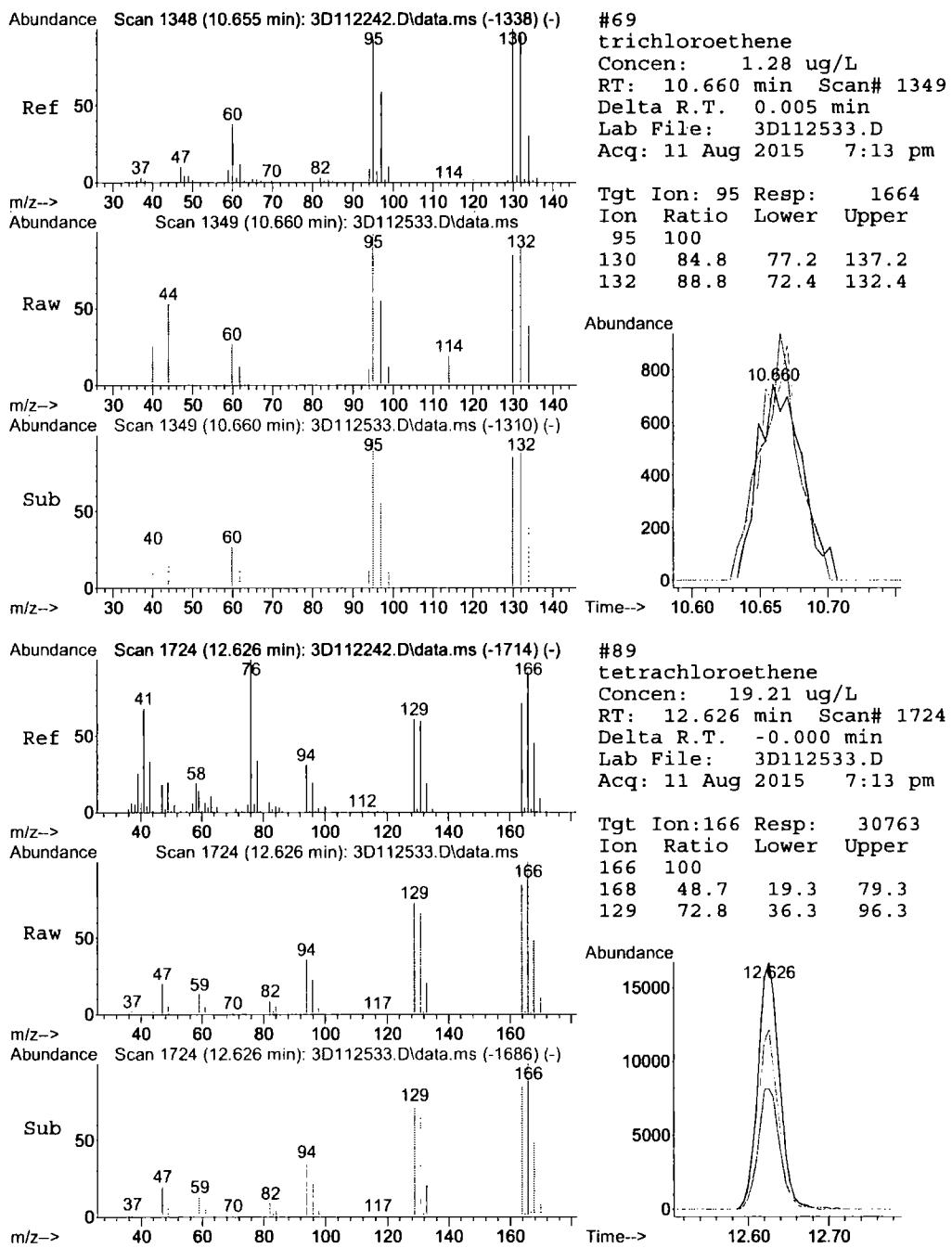
Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112533.D
 Acq On : 11 Aug 2015 7:13 pm
 Operator : ximenac
 Sample : jc1106-15
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 26 Sample Multiplier: 1

Quant Time: Aug 12 09:29:30 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration









Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\

Data File : 3D112573.D

Acq On : 12 Aug 2015 1:54 pm

Operator : ximenac

Sample : jc1106-16

Misc : MS89468,V3D4824,5,,,1

ALS Vial : 12 Sample Multiplier: 1

Quant Time: Aug 12 14:32:28 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

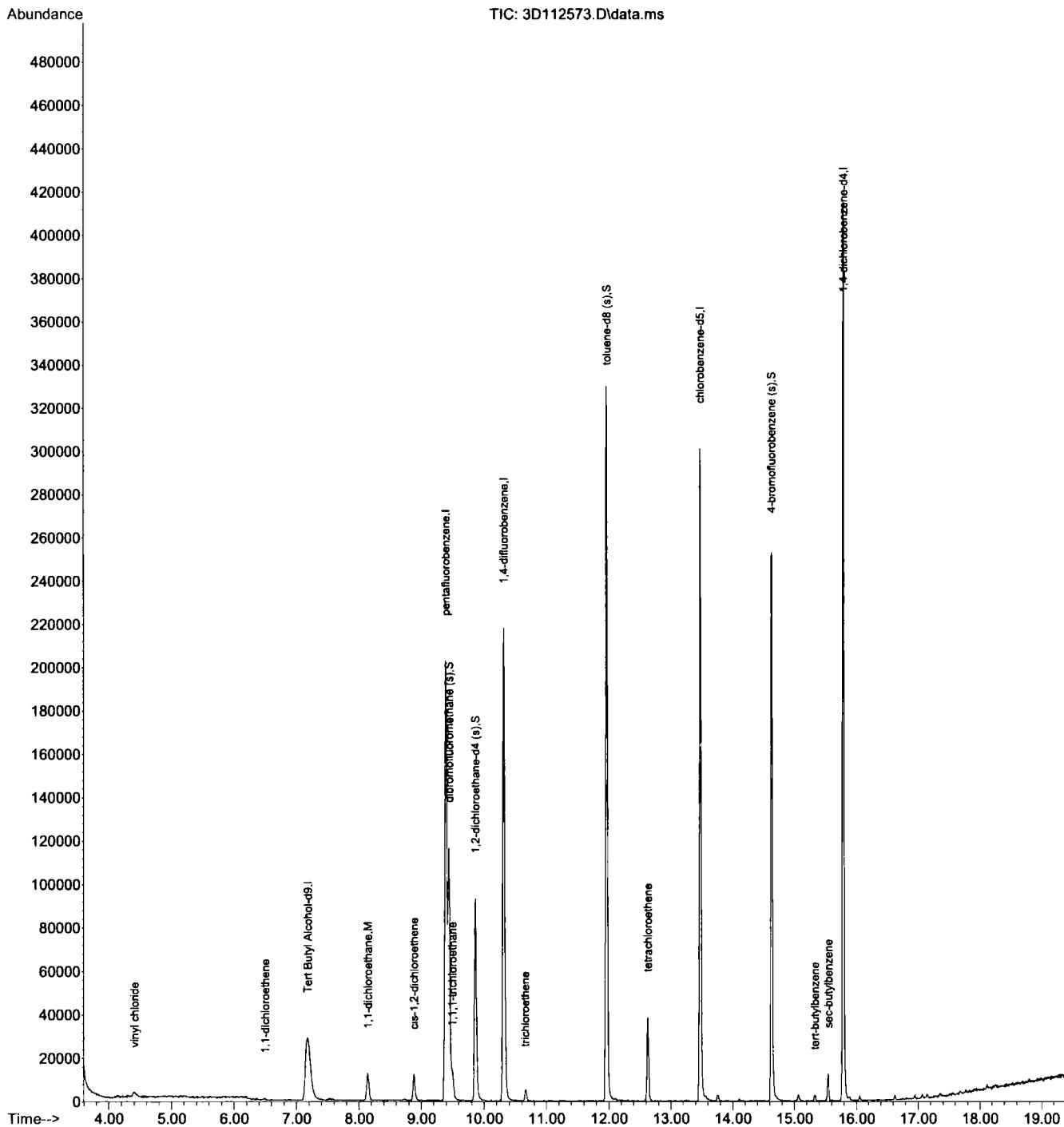
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.178	65	88764	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	186964	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.313	114	217383	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	181788	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	113725	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	79219	55.30	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	110.60%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	74462	56.11	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	112.22%	
80) toluene-d8 (s)	11.965	98	238139	50.02	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	100.04%	
106) 4-bromofluorobenzene (s)	14.628	95	83477	47.71	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	95.42%	
Target Compounds						
11) vinyl chloride	4.405	62	4291	1.58	ug/L	90
22) 1,1-dichloroethene	6.497	61	475	0.21	ug/L	# 60
37) 1,1-dichloroethane	8.143	63	16610	6.89	ug/L	97
43) cis-1,2-dichloroethene	8.882	96	7003	4.30	ug/L	# 80
54) 1,1,1-trichloroethane	9.496	97	10019	4.17	ug/L	96
69) trichloroethene	10.654	95	2269	1.84	ug/L	84
89) tetrachloroethene	12.626	166	13672	8.83	ug/L	92
117) tert-butylbenzene	15.331	119	1604	0.40	ug/L	80
120) sec-butylbenzene	15.546	105	8320	1.30	ug/L	91

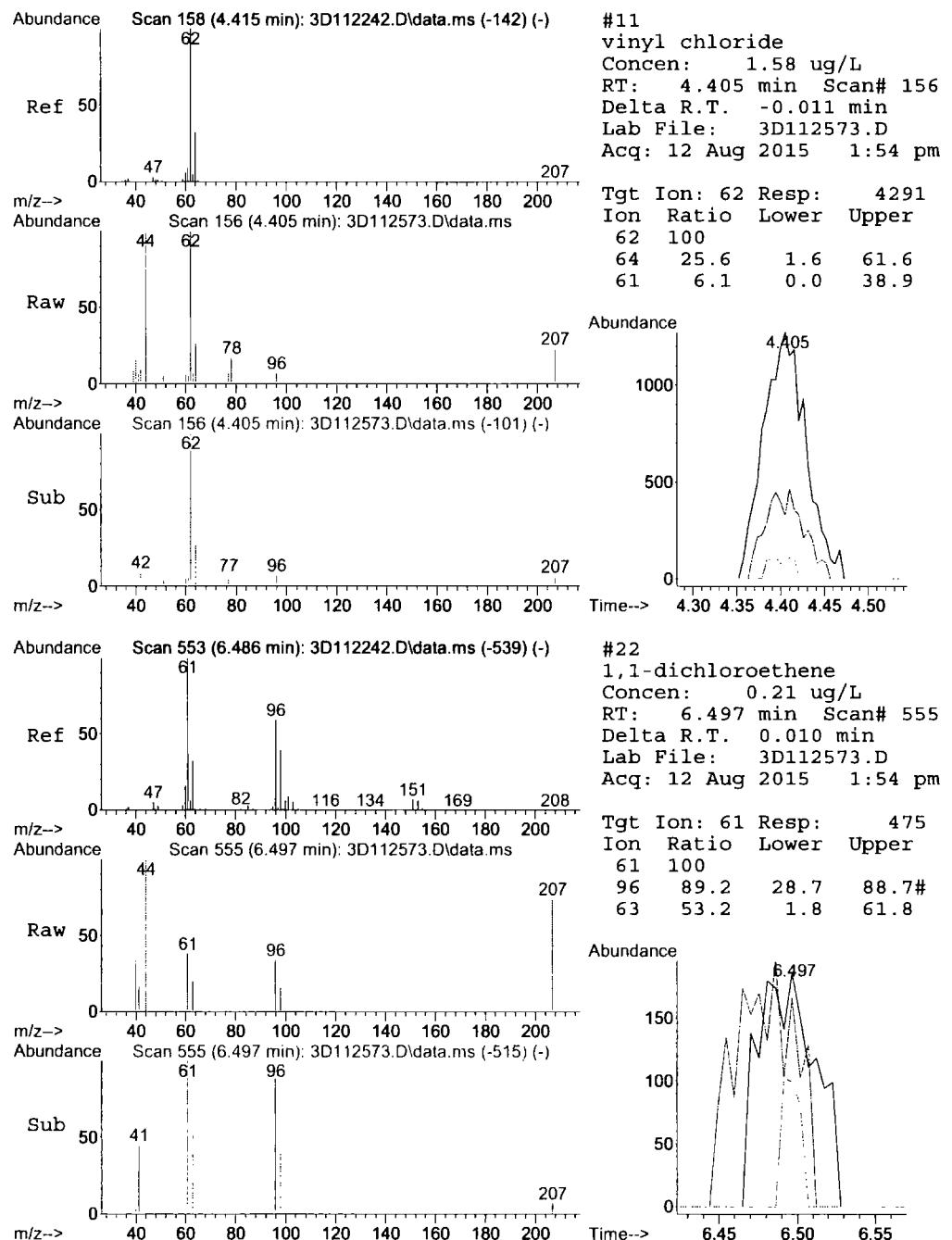
(#= qualifier out of range (m) = manual integration (+) = signals summed

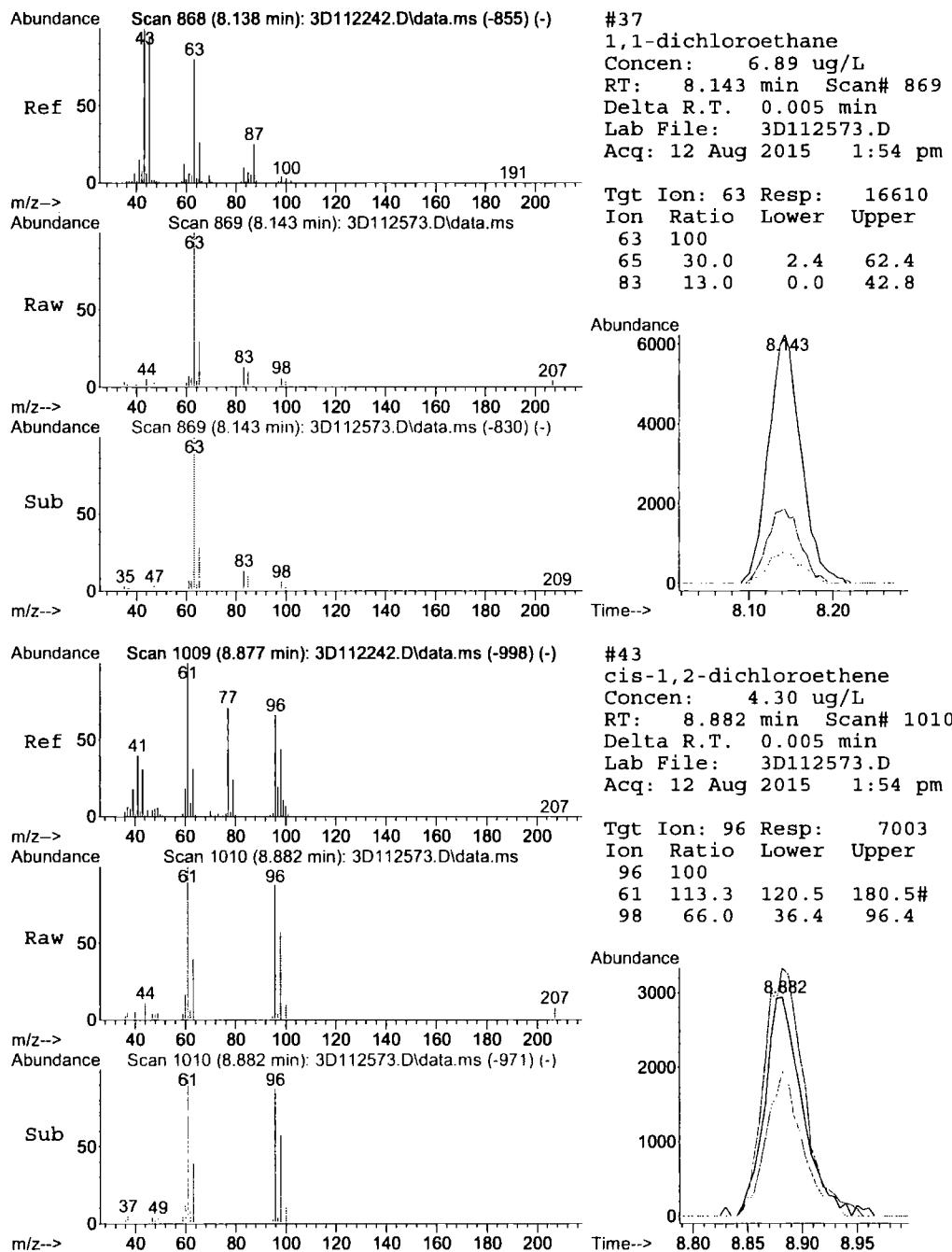
Quantitation Report (QT Reviewed)

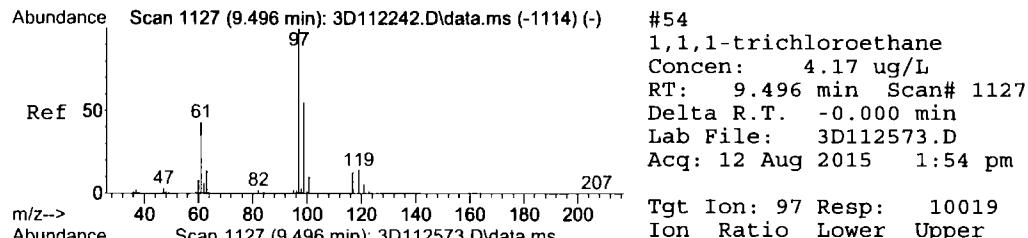
Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
 Data File : 3D112573.D
 Acq On : 12 Aug 2015 1:54 pm
 Operator : ximenac
 Sample : jc1106-16
 Misc : MS89468,V3D4824,5,,,1
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Aug 12 14:32:28 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

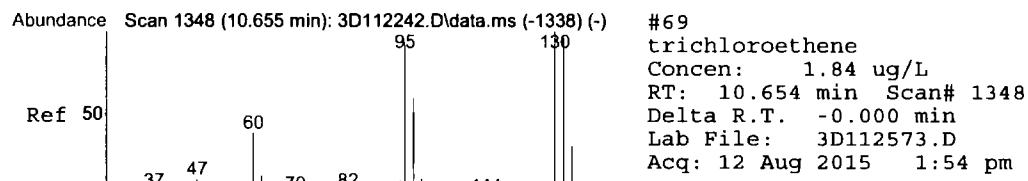
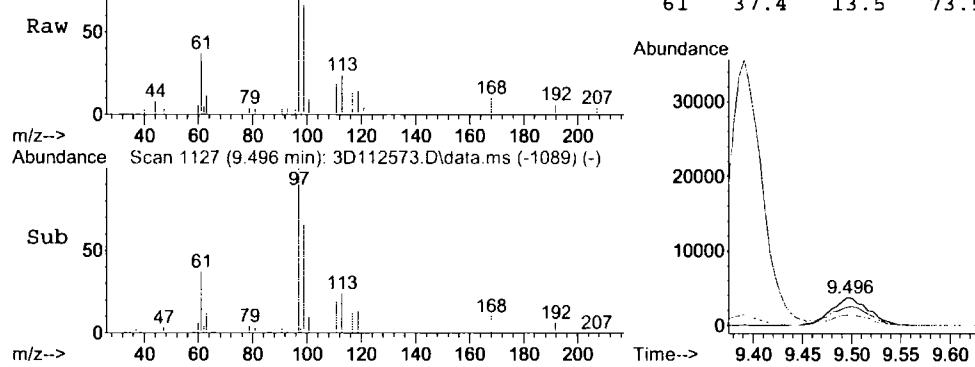




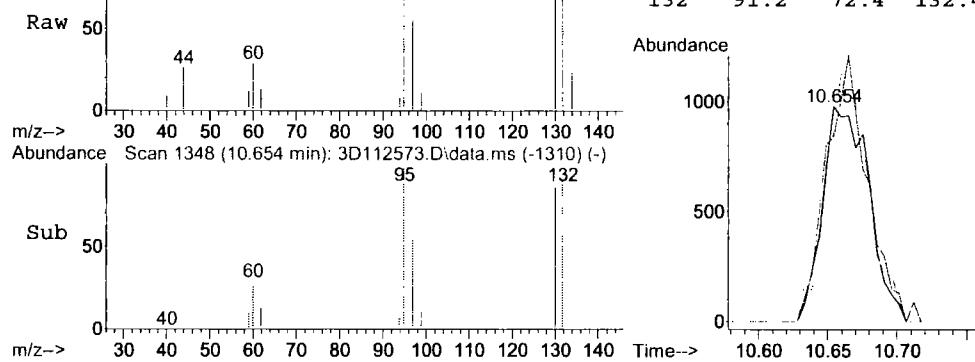


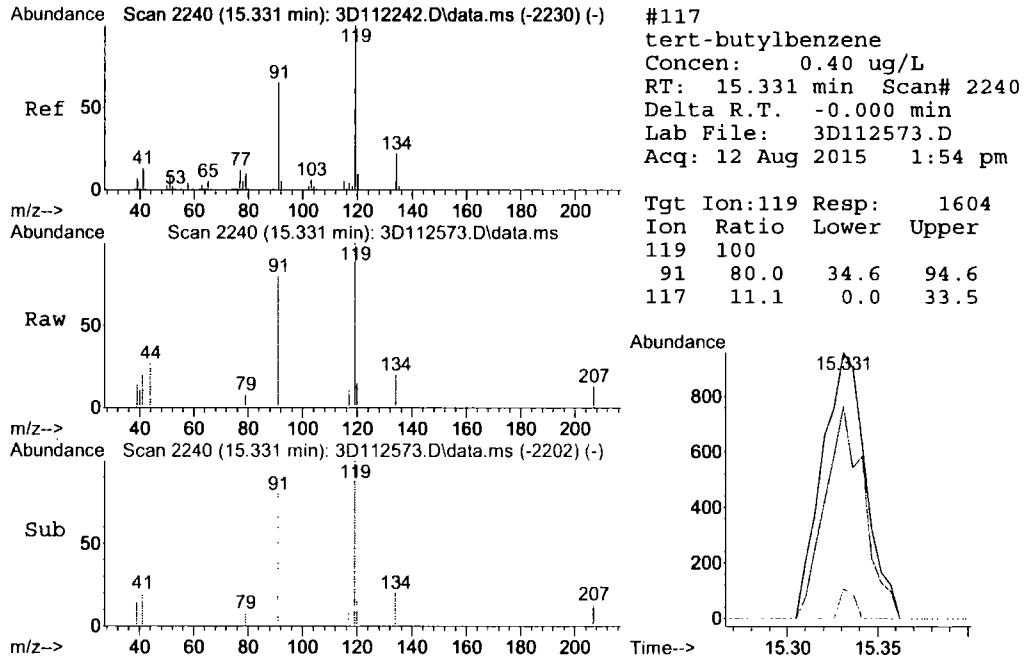
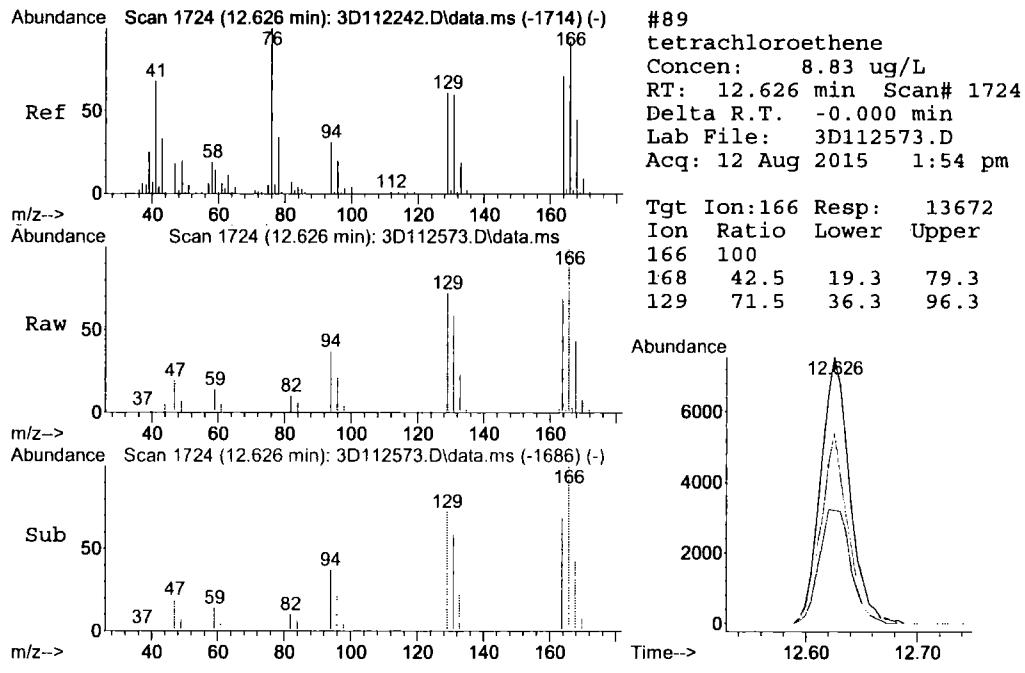


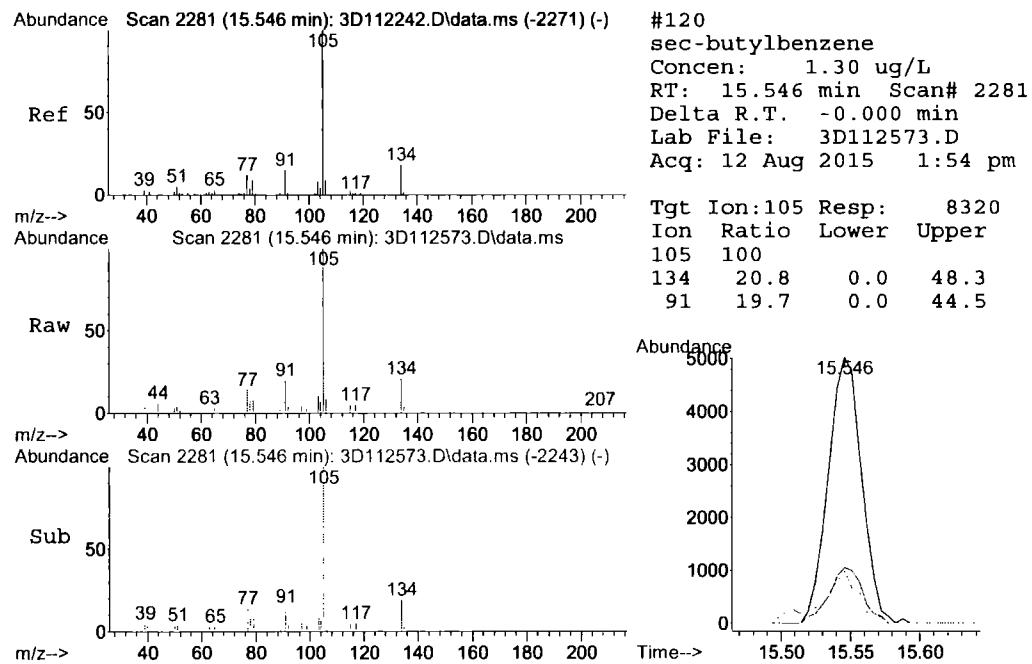
Tgt Ion:	97	Ion Ratio	100	Resp:	10019
Ion	97	100		Lower	
	99	63.7	33.7	Upper	93.7
	61	37.4	13.5		73.5



Tgt Ion:	95	Ion Ratio	100	Resp:	2269
Ion	95	100		Lower	
	130	86.3	77.2	Upper	137.2
	132	91.2	72.4		132.4







7.1.16

7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112555.D
 Acq On : 12 Aug 2015 5:07 am
 Operator : ximenac
 Sample : jc1106-17
 Misc : MS89468,V3D4823,5,,,1
 ALS Vial : 48 Sample Multiplier: 1

Quant Time: Aug 12 09:45:45 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

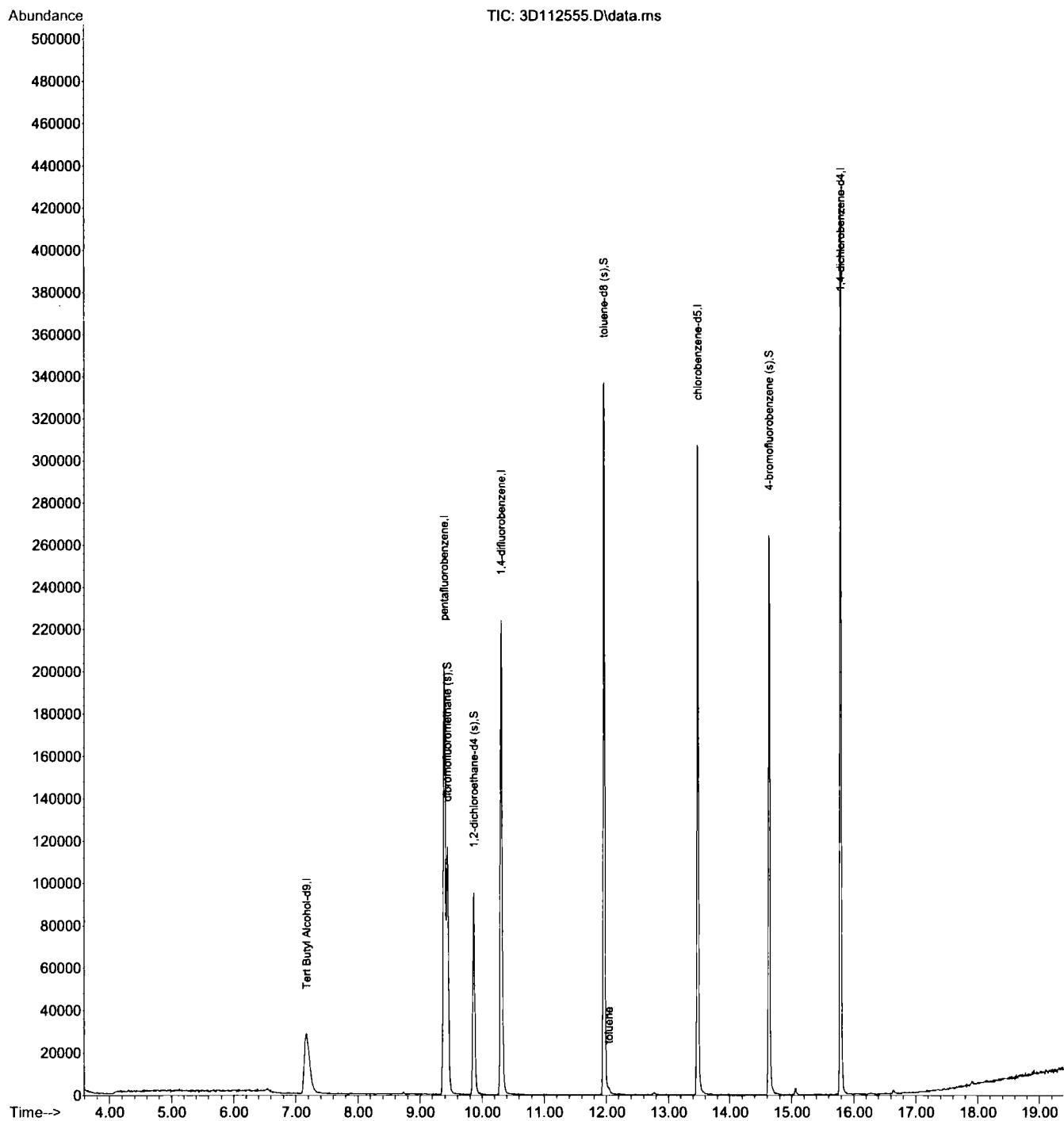
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	94319	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	190266	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	224022	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	184821	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	114314	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.438	113	78516	53.86	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	=	107.72%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	75635	56.01	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	=	112.02%	
80) toluene-d8 (s)	11.960	98	247330	50.41	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	=	100.82%	
106) 4-bromofluorobenzene (s)	14.629	95	86055	48.93	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	=	97.86%	
Target Compounds						
82) toluene	12.039	92	965	0.33	ug/L	# 79

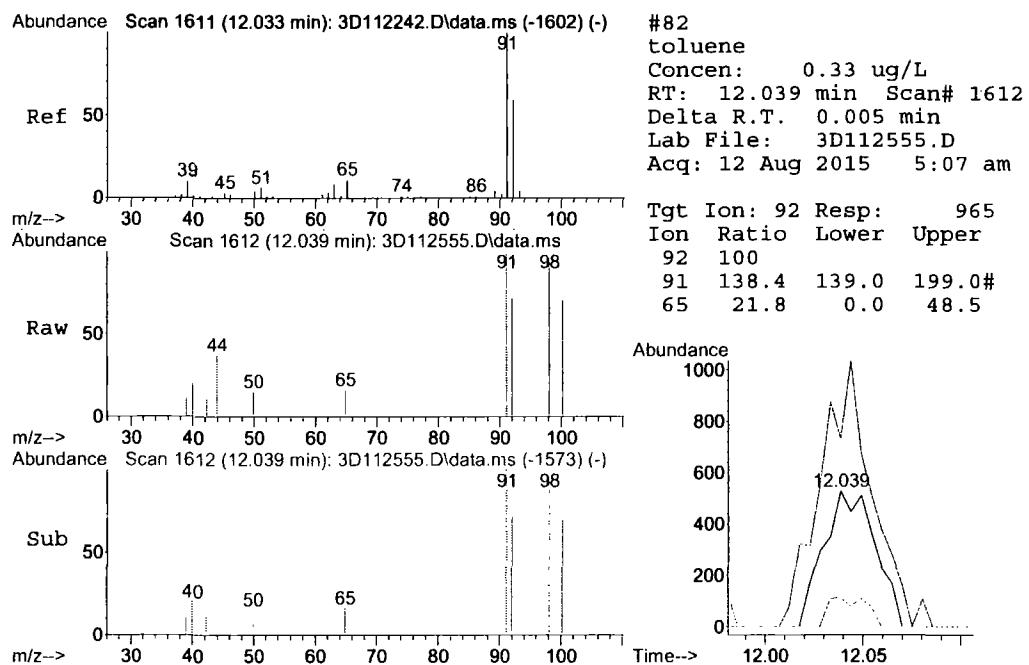
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112555.D
 Acq On : 12 Aug 2015 5:07 am
 Operator : ximenac
 Sample : jc1106-17
 Misc : MS89468,V3D4823,5,,,1
 ALS Vial : 48 Sample Multiplier: 1

Quant Time: Aug 12 09:45:45 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\

Data File : 3D112534.D

Acq On : 11 Aug 2015 7:40 pm

Operator : ximenac

Sample : jc1106-18

Misc : MS89468,V3D4822,5,,,1

ALS Vial : 27 Sample Multiplier: 1

Quant Time: Aug 12 09:30:27 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

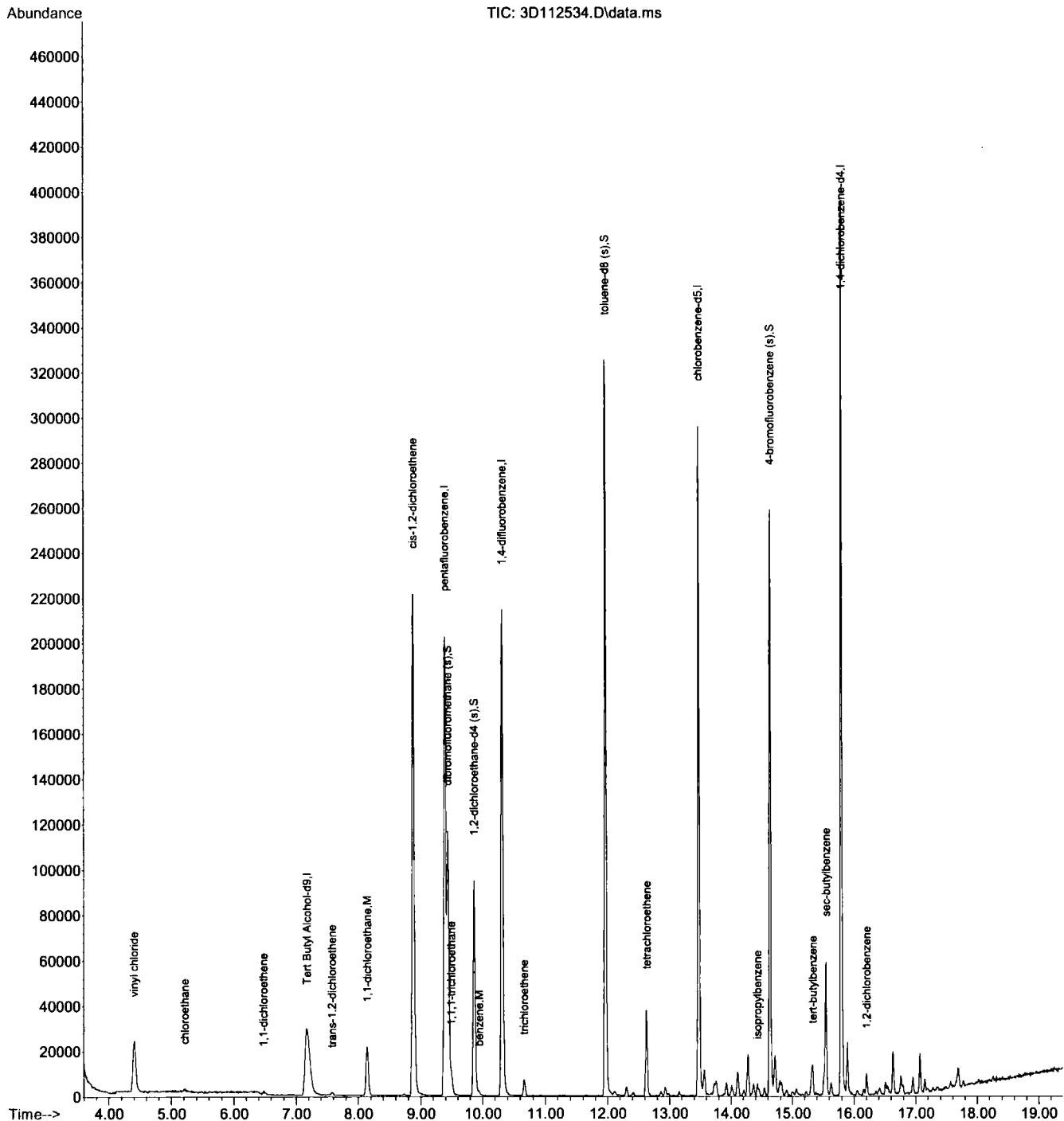
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	88719	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	186366	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	213690	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	175269	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.793	152	108187	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.438	113	79434	55.63	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	=	111.26%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	73921	55.89	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	=	111.78%	
80) toluene-d8 (s)	11.960	98	235002	50.21	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	=	100.42%	
106) 4-bromofluorobenzene (s)	14.634	95	81667	49.06	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	=	98.12%	
Target Compounds						
					Qvalue	
11) vinyl chloride	4.399	62	44111	16.33	ug/L	98
13) chloroethane	5.212	64	1656	1.45	ug/L	86
22) 1,1-dichloroethene	6.476	61	1088	0.49	ug/L	90
33) trans-1,2-dichloroethene	7.577	61	1028	0.51	ug/L	82
37) 1,1-dichloroethane	8.143	63	29062	12.09	ug/L	99
43) cis-1,2-dichloroethene	8.877	96	122761	75.58	ug/L	89
54) 1,1,1-trichloroethane	9.496	97	7363	3.08	ug/L	95
65) benzene	9.941	78	1730	0.34	ug/L	89
69) trichloroethene	10.660	95	2852	2.35	ug/L	91
89) tetrachloroethene	12.631	166	13486	9.04	ug/L	95
105) isopropylbenzene	14.429	105	3775	0.70	ug/L	94
117) tert-butylbenzene	15.336	119	6012	1.58	ug/L	99
120) sec-butylbenzene	15.546	105	42171	6.93	ug/L	98
124) 1,2-dichlorobenzene	16.202	146	4175	1.30	ug/L	95

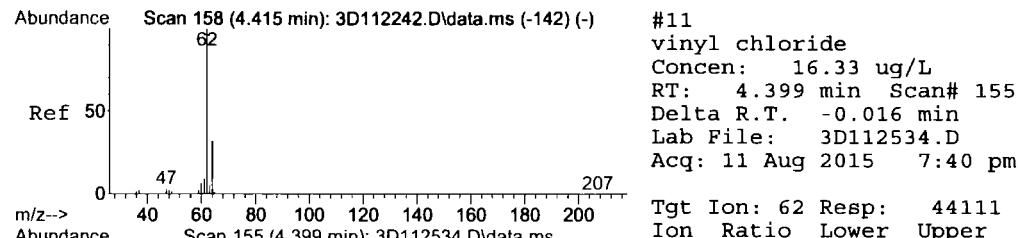
(#= qualifier out of range (m)= manual integration (+)= signals summed

Quantitation Report (QT Reviewed)

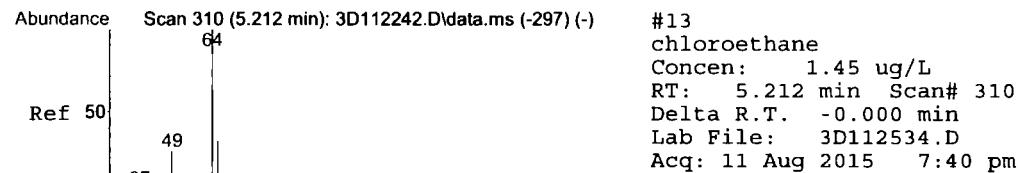
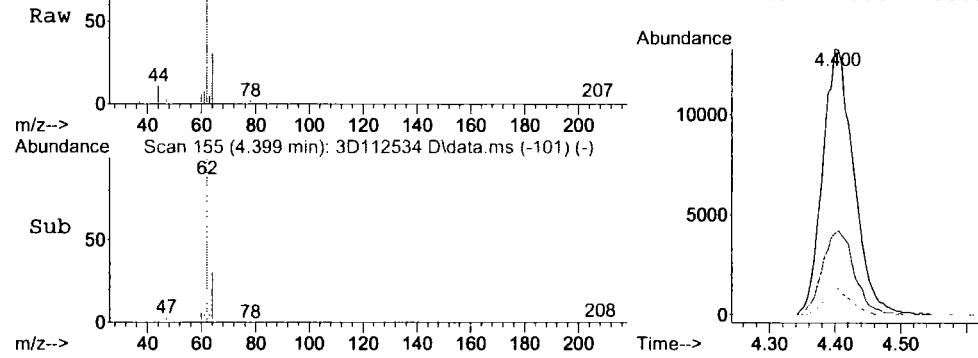
Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112534.D
 Acq On : 11 Aug 2015 7:40 pm
 Operator : ximenac
 Sample : jc1106-18
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Aug 12 09:30:27 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

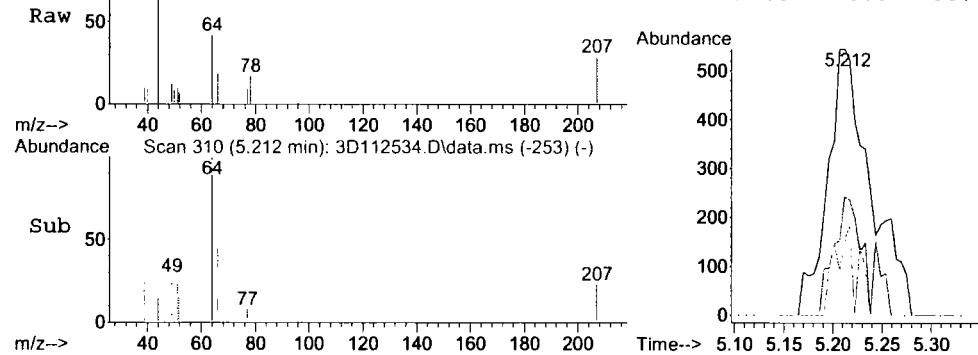


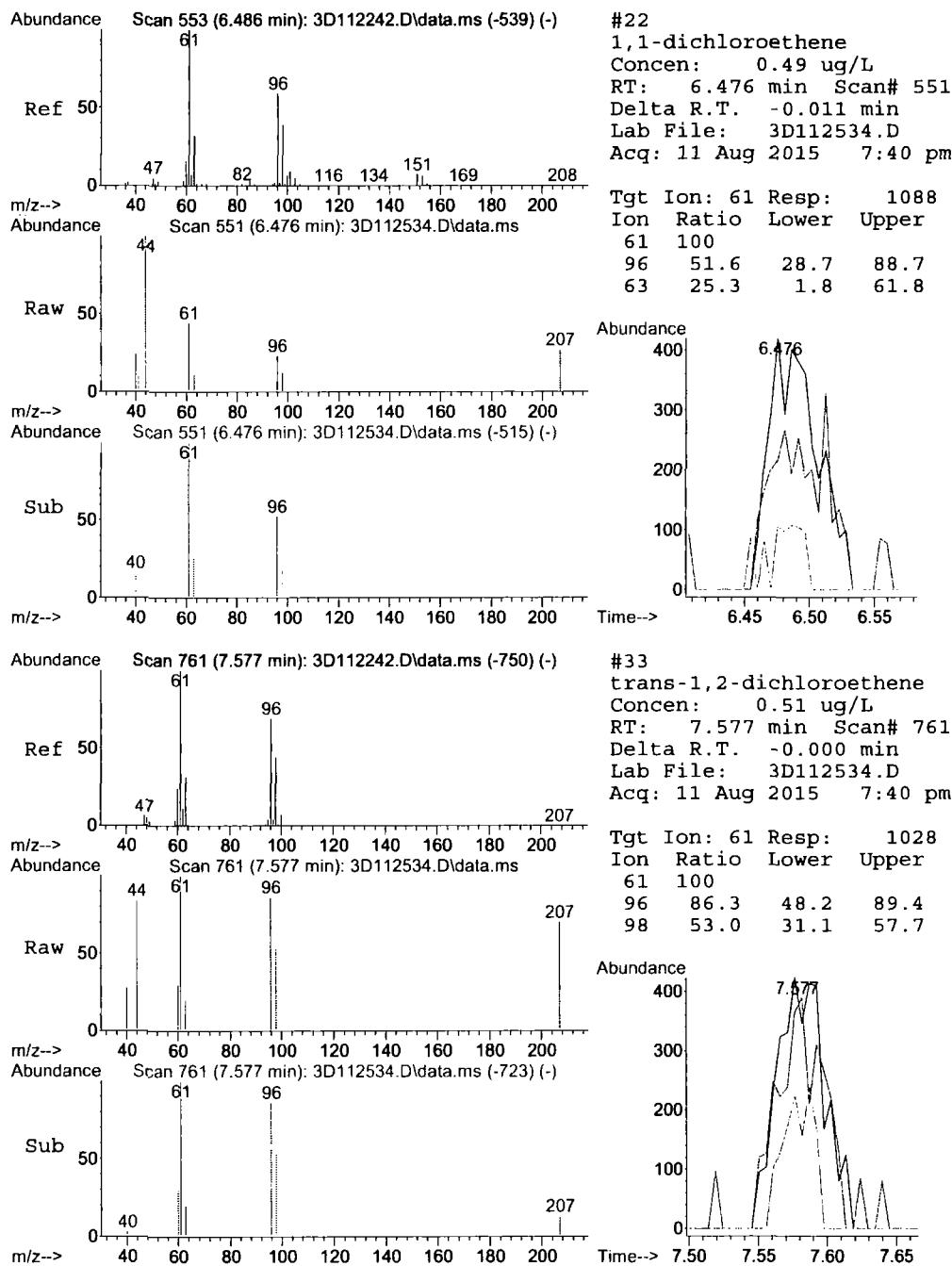


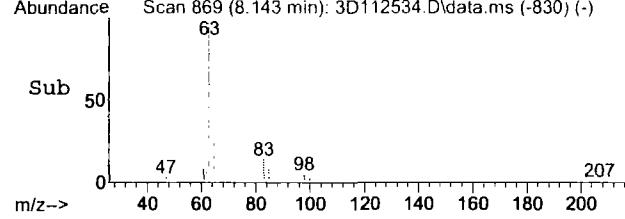
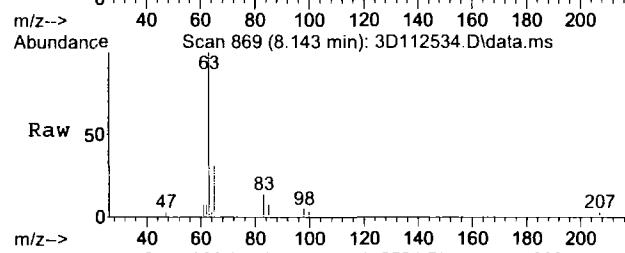
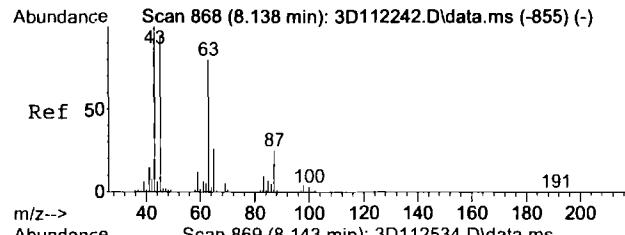
Tgt	Ion:	62	Resp:	44111
Ion	Ratio	100		
62				
64	30.6	1.6	61.6	
61	7.9	0.0	38.9	



Tgt	Ion:	64	Resp:	1656
Ion	Ratio	100		
64				
66	44.6	1.6	61.6	
49	27.5	0.0	56.0	

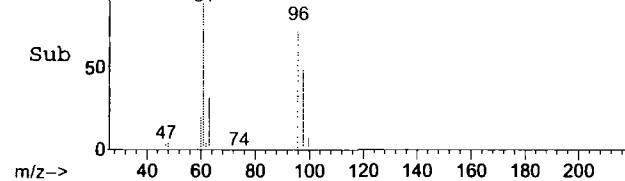
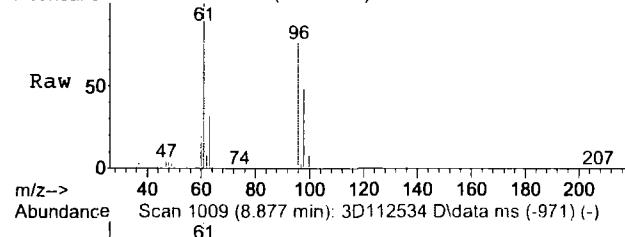
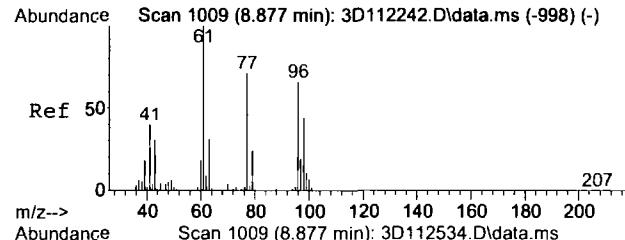
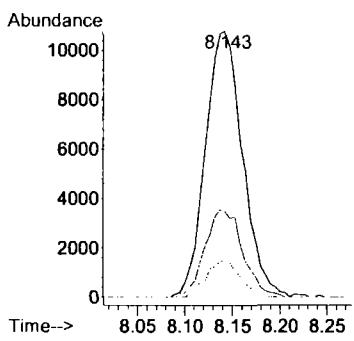






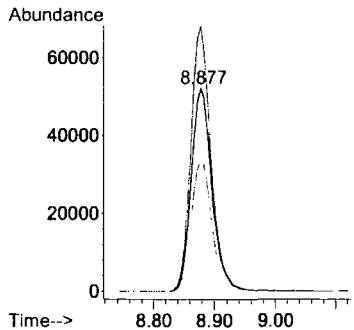
#37
1,1-dichloroethane
Concen: 12.09 ug/L
RT: 8.143 min Scan# 869
Delta R.T. 0.005 min
Lab File: 3D112534.D
Acq: 11 Aug 2015 7:40 pm

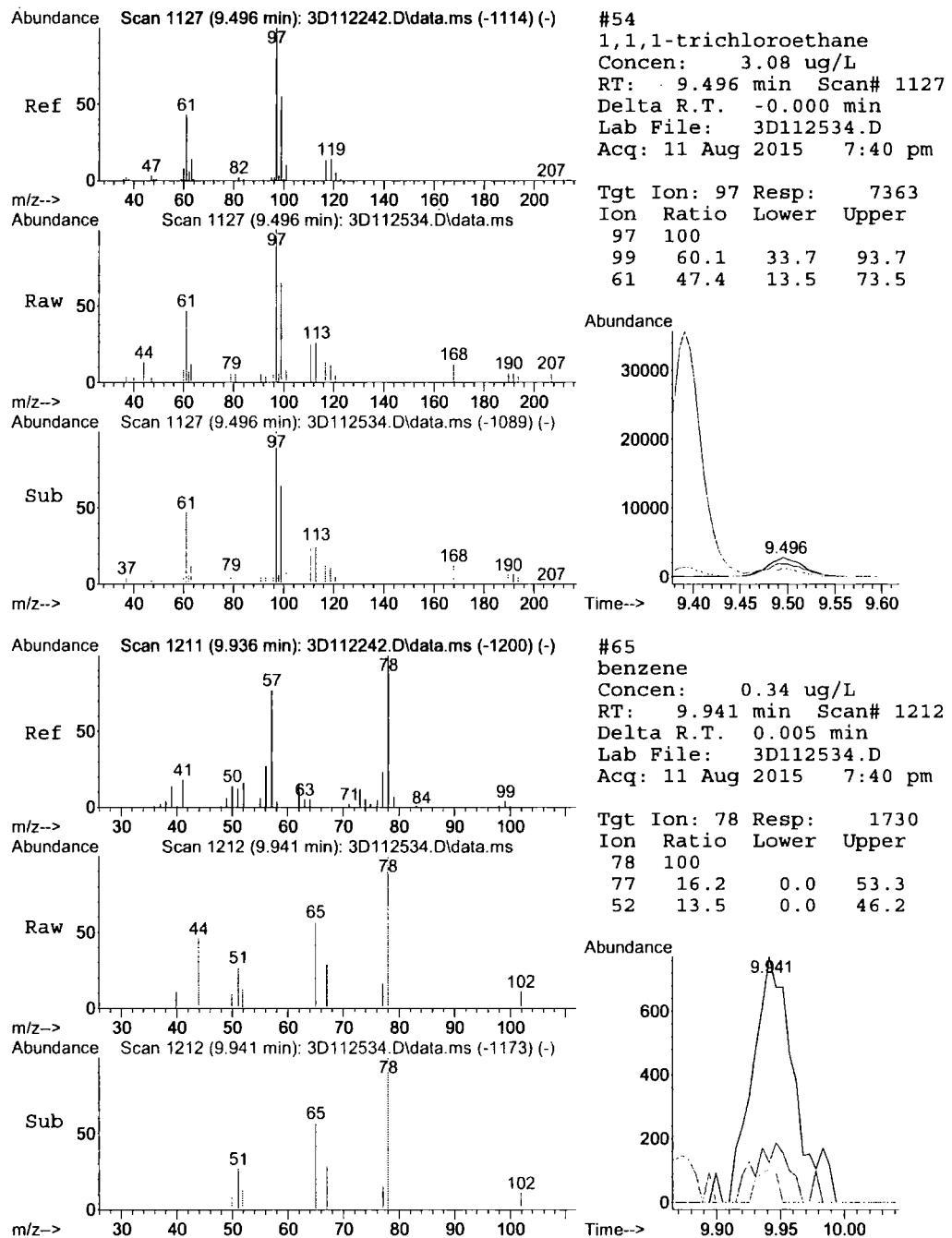
Tgt	Ion:	63	Resp:	29062
Ion	Ratio	100	Lower	
63	100			
65	32.3	2.4	62.4	
83	13.8	0.0	42.8	

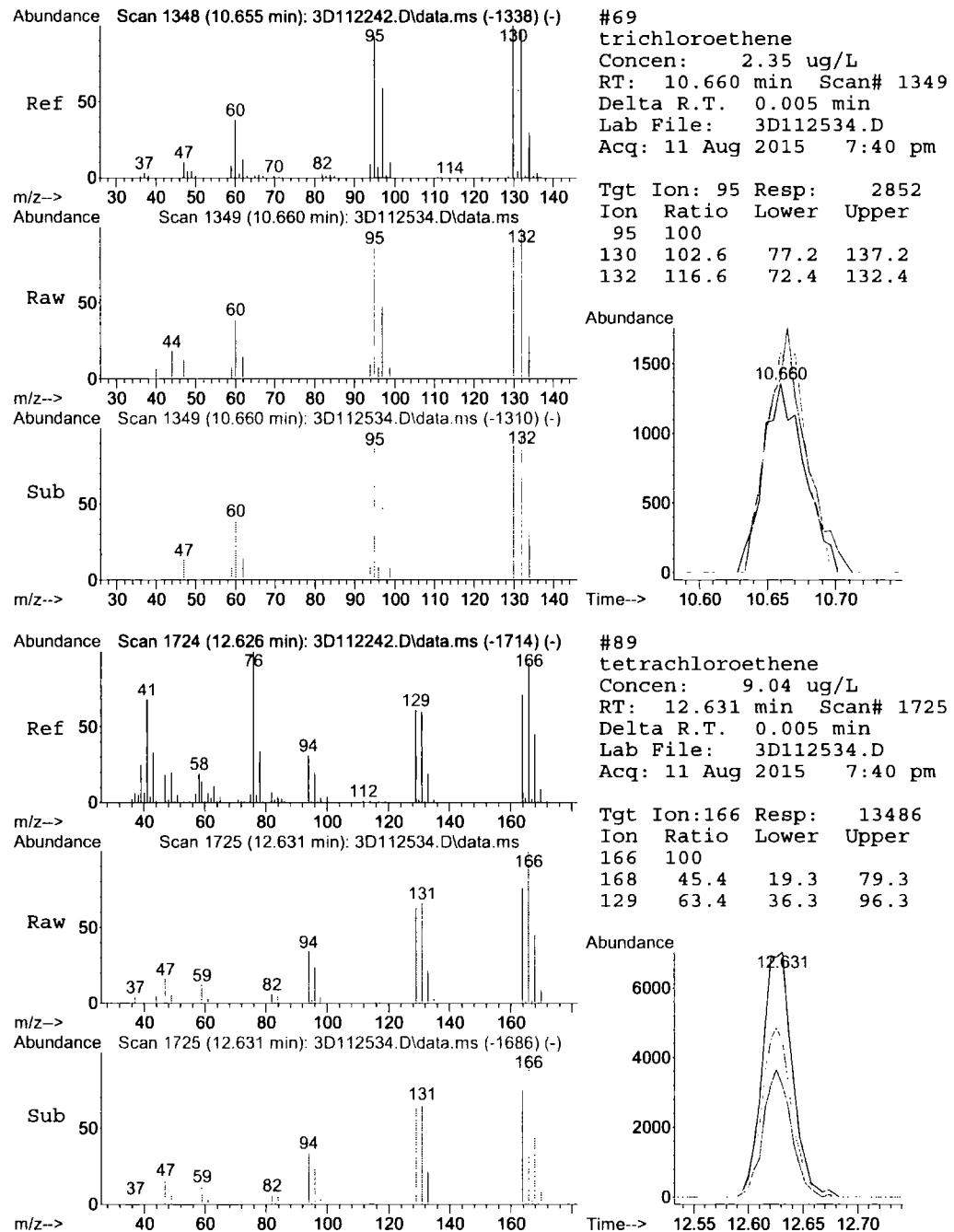


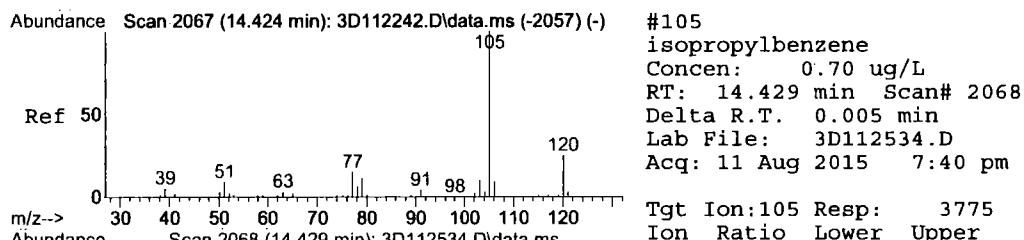
#43
cis-1,2-dichloroethene
Concen: 75.58 ug/L
RT: 8.877 min Scan# 1009
Delta R.T. -0.000 min
Lab File: 3D112534.D
Acq: 11 Aug 2015 7:40 pm

Tgt	Ion:	96	Resp:	122761
Ion	Ratio	100	Lower	
96	100			
61	130.9	120.5	180.5	
98	64.3	36.4	96.4	

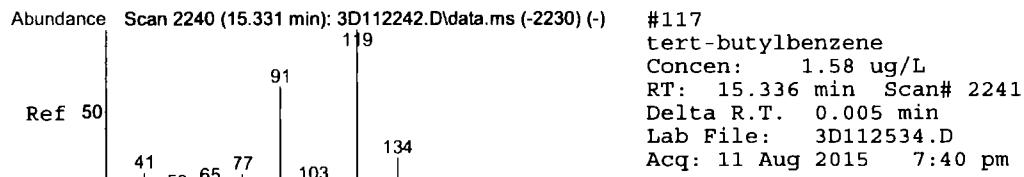
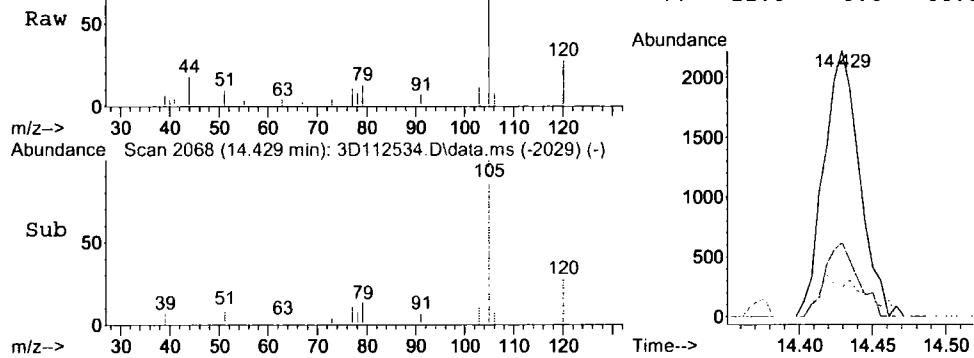




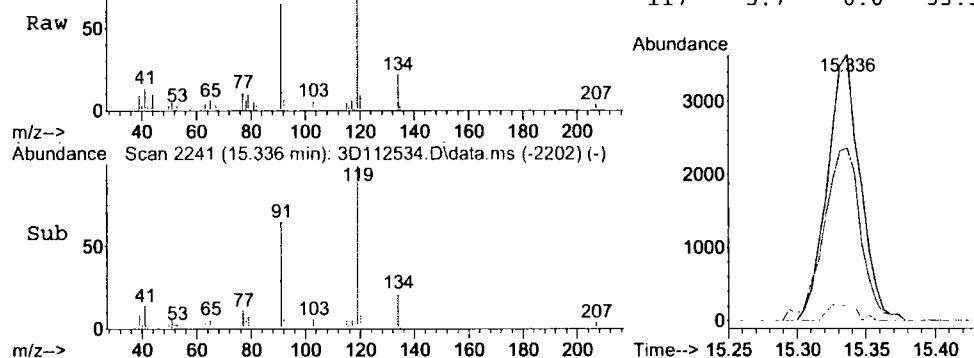


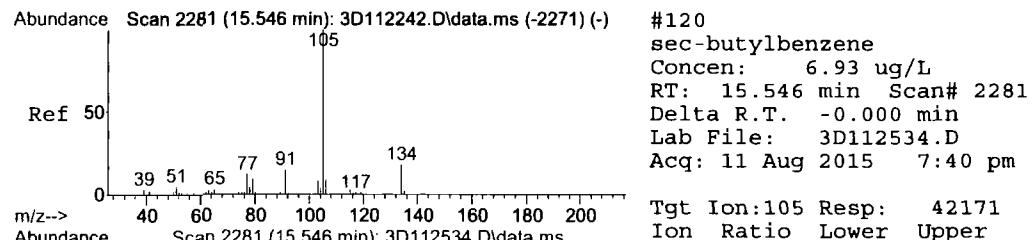


Tgt Ion:105 Resp: 3775
Ion Ratio Lower Upper
105 100
120 27.8 0.0 55.4
77 11.0 0.0 44.8

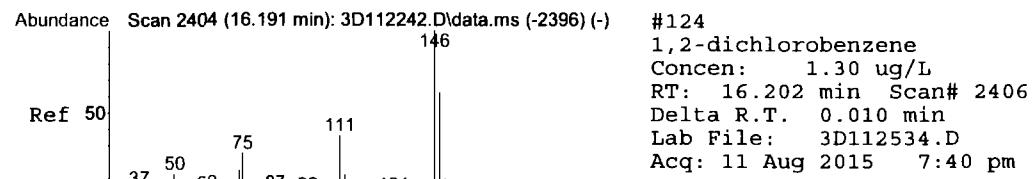
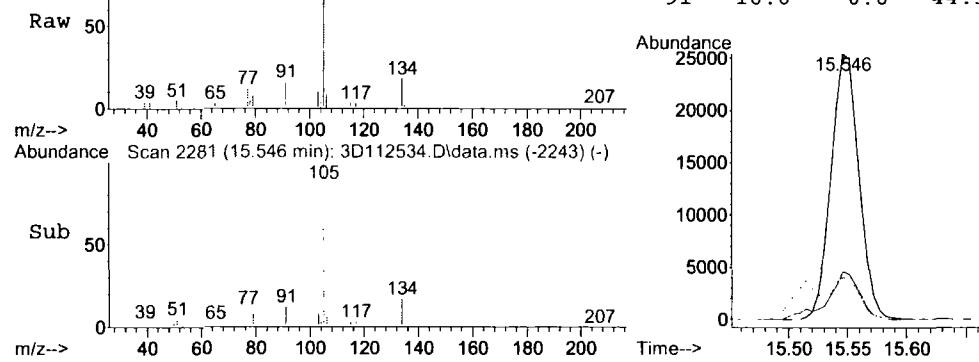


Tgt Ion:119 Resp: 6012
Ion Ratio Lower Upper
119 100
91 64.9 34.6 94.6
117 5.7 0.0 33.5

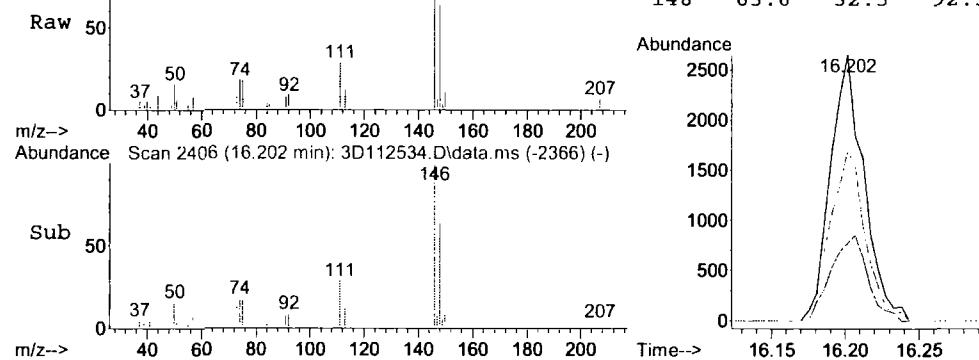




Tgt	Ion:105	Resp:	42171
105	100		
134	18.1	0.0	48.3
91	16.0	0.0	44.5



Tgt	Ion:146	Resp:	4175
146	100		
111	29.1	6.4	66.4
148	63.6	32.5	92.5



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112529.D
 Acq On : 11 Aug 2015 5:24 pm
 Operator : ximenac
 Sample : jc1106-19
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Aug 12 09:28:12 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

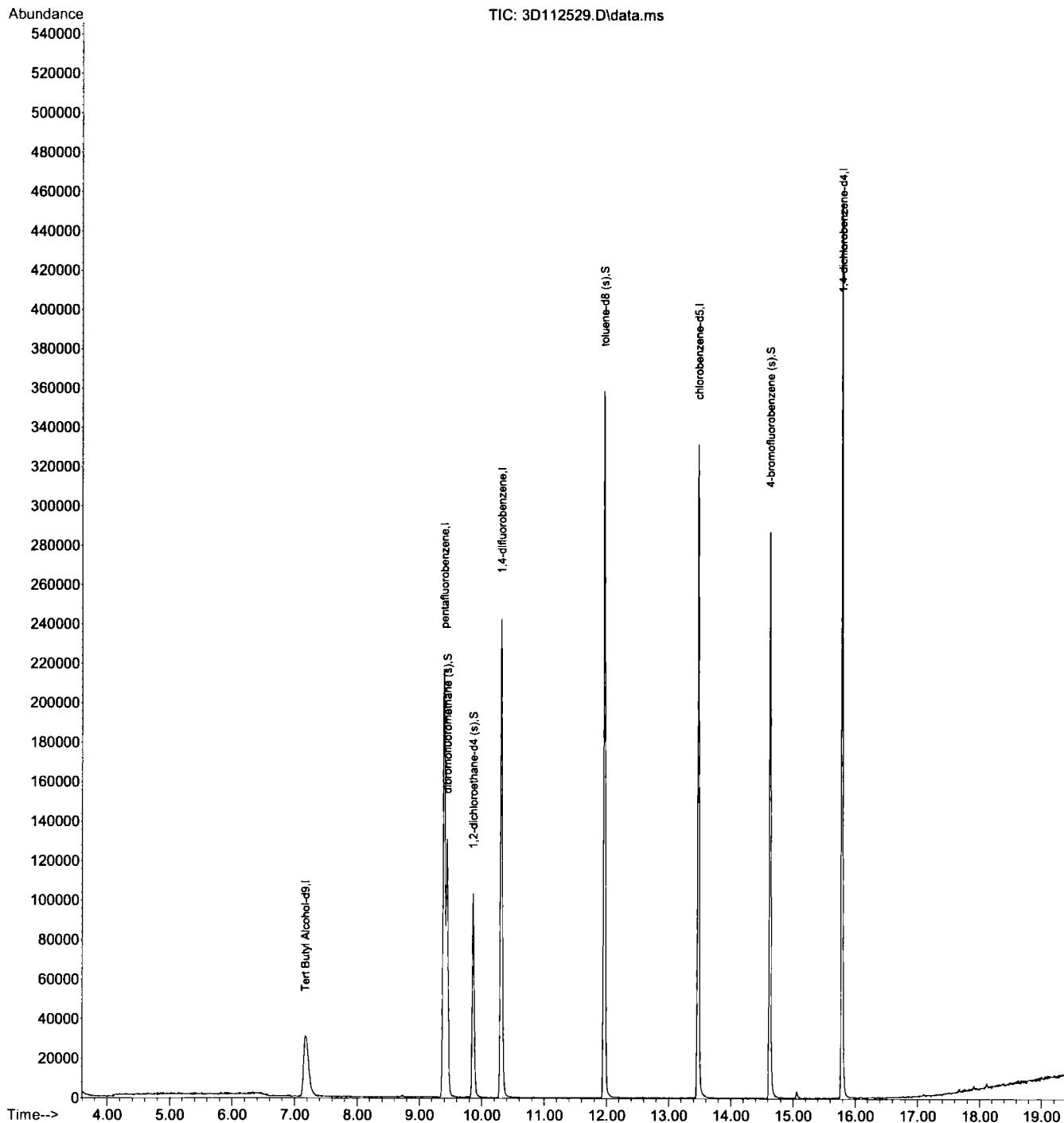
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.168	65	97633	500.00	ug/L	-0.01
4) pentafluorobenzene	9.396	168	205375	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	241116	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	197855	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	123858	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	88077	55.97	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	111.94%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	82016	56.27	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	112.54%	
80) toluene-d8 (s)	11.960	98	265412	50.26	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	100.52%	
106) 4-bromofluorobenzene (s)	14.629	95	91661	48.10	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	96.20%	

Target Compounds	Qvalue
(#= qualifier out of range (m)= manual integration (+)= signals summed	

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112529.D
 Acq On : 11 Aug 2015 5:24 pm
 Operator : ximenac
 Sample : jc1106-19
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Aug 12 09:28:12 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112519.D
 Acq On : 11 Aug 2015 12:38 pm
 Operator : ximenac
 Sample : mb
 Misc : MS88280,V3D4822,5,,,1
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Aug 11 16:10:33 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.178	65	103952	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	226869	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	264986	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	216656	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.793	152	129511	50.00	ug/L	0.00

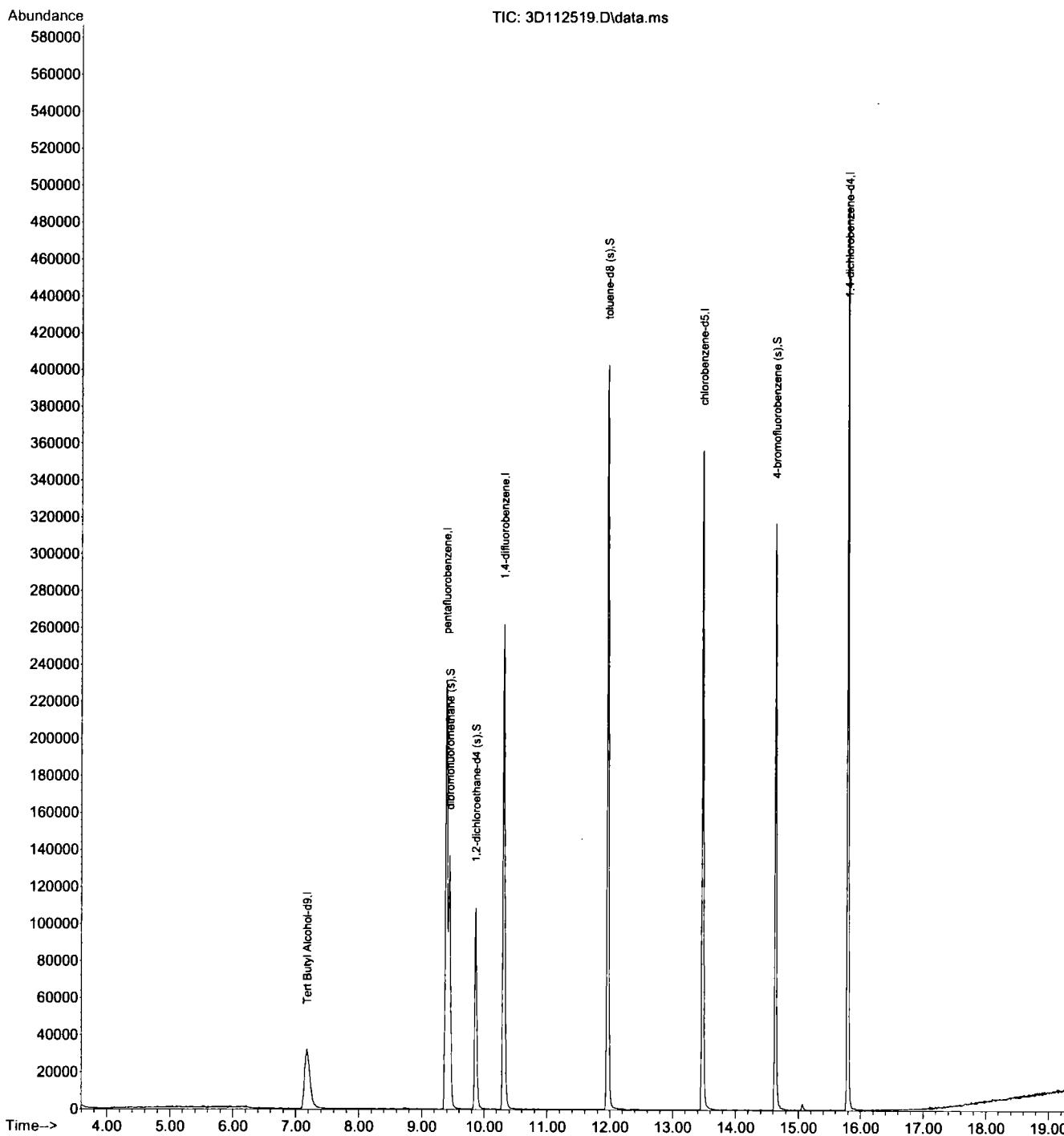
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	93639	53.87	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	107.74%
51) 1,2-dichloroethane-d4 (s)	9.863	65	89940	55.86	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	111.72%
80) toluene-d8 (s)	11.965	98	294612	50.76	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	101.52%
106) 4-bromofluorobenzene (s)	14.634	95	99709	50.04	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	100.08%

Target Compounds	Qvalue
(#) = qualifier out of range (m) = manual integration (+) = signals summed	

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112519.D
 Acq On : 11 Aug 2015 12:38 pm
 Operator : ximenac
 Sample : mb
 Misc : MS88280,V3D4822,5,,,1
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Aug 11 16:10:33 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112544.D
 Acq On : 12 Aug 2015 12:11 am
 Operator : ximenac
 Sample : mb
 Misc : MS89457,V3D4823,5,,,1
 ALS Vial : 37 Sample Multiplier: 1

Quant Time: Aug 12 09:40:24 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.163	65	82447	500.00	ug/L	-0.02
4) pentafluorobenzene	9.391	168	171750	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	198901	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	165044	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	101680	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	73569	55.91	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	111.82%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	69179	56.75	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	113.50%	
80) toluene-d8 (s)	11.960	98	222103	50.99	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	101.98%	
106) 4-bromofluorobenzene (s)	14.629	95	75250	48.10	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	96.20%	

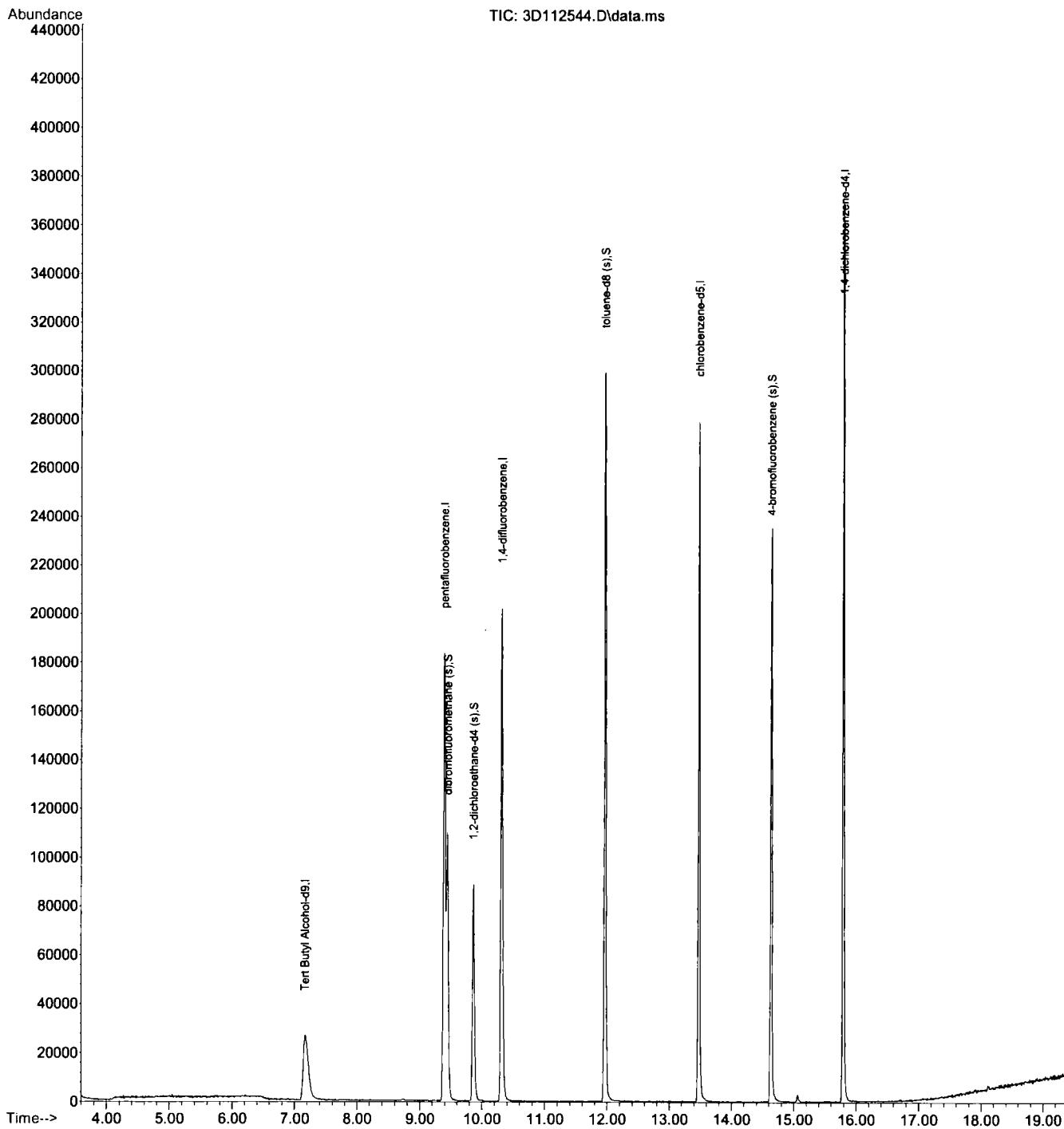
Target Compounds	Qvalue
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112544.D
 Acq On : 12 Aug 2015 12:11 am
 Operator : ximenac
 Sample : mb
 Misc : MS89457,V3D4823,5,,,1
 ALS Vial : 37 Sample Multiplier: 1

Quant Time: Aug 12 09:40:24 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112569.D
 Acq On : 12 Aug 2015 12:06 pm
 Operator : ximenac
 Sample : mb
 Misc : MS89457, V3D4824, 5, , , 1
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Aug 12 14:30:59 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.184	65	95029	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	192060	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	221163	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	183575	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	114634	50.00	ug/L	0.00

System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	83863	56.99	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	113.98%
51) 1,2-dichloroethane-d4 (s)	9.863	65	75047	55.05	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	110.10%
80) toluene-d8 (s)	11.965	98	245477	50.68	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	101.36%
106) 4-bromofluorobenzene (s)	14.629	95	85262	48.34	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.68%

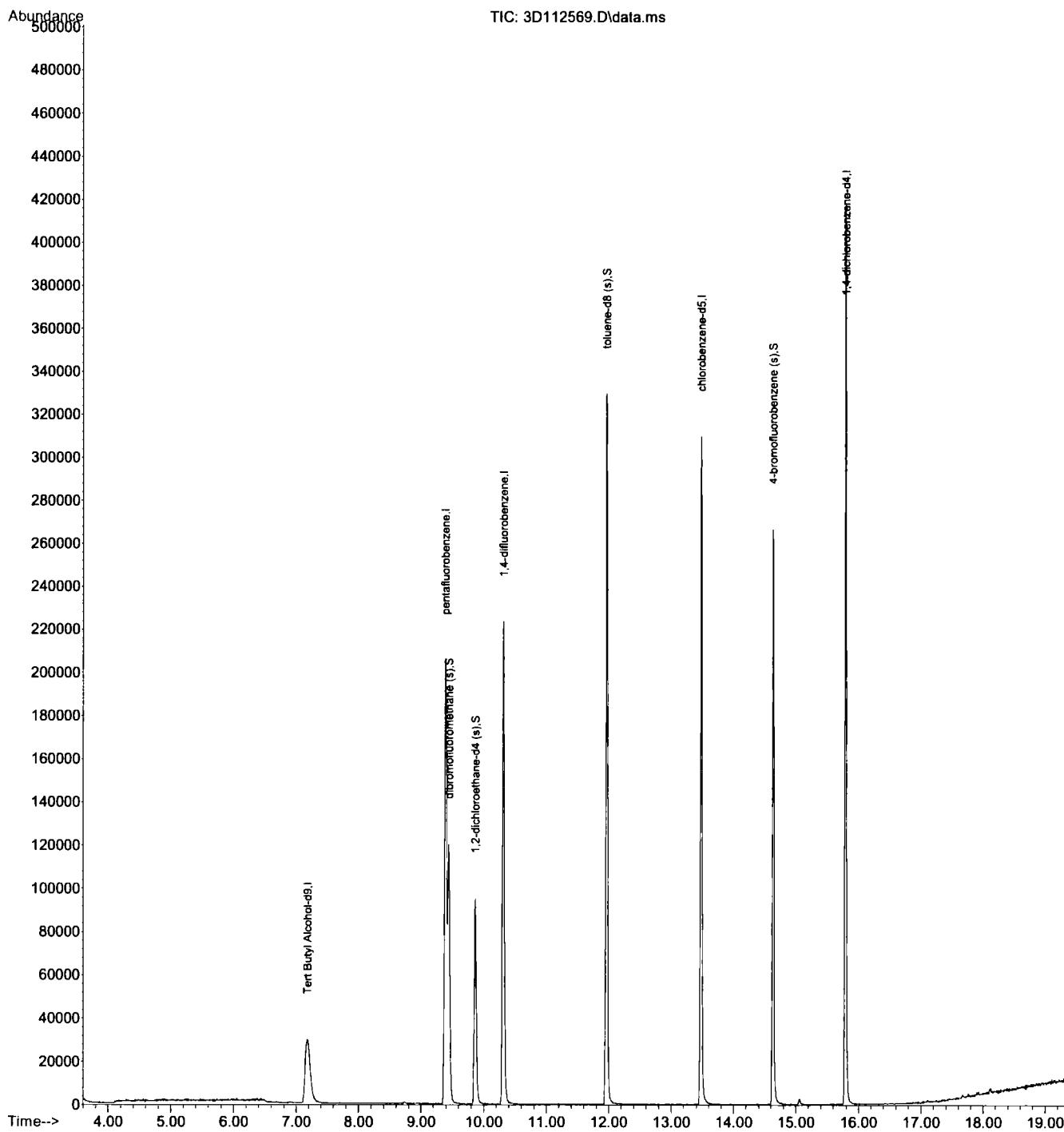
Target Compounds	Qvalue
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112569.D
 Acq On : 12 Aug 2015 12:06 pm
 Operator : ximenac
 Sample : mb
 Misc : MS89457,V3D4824,5,,,1
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Aug 12 14:30:59 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112565.D
 Acq On : 12 Aug 2015 10:15 am
 Operator : ximenac
 Sample : mb2
 Misc : MS89457,V3D4823,5,,,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Aug 12 14:29:31 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.184	65	89468	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	182832	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	209972	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	174088	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	109833	50.00	ug/L	0.00

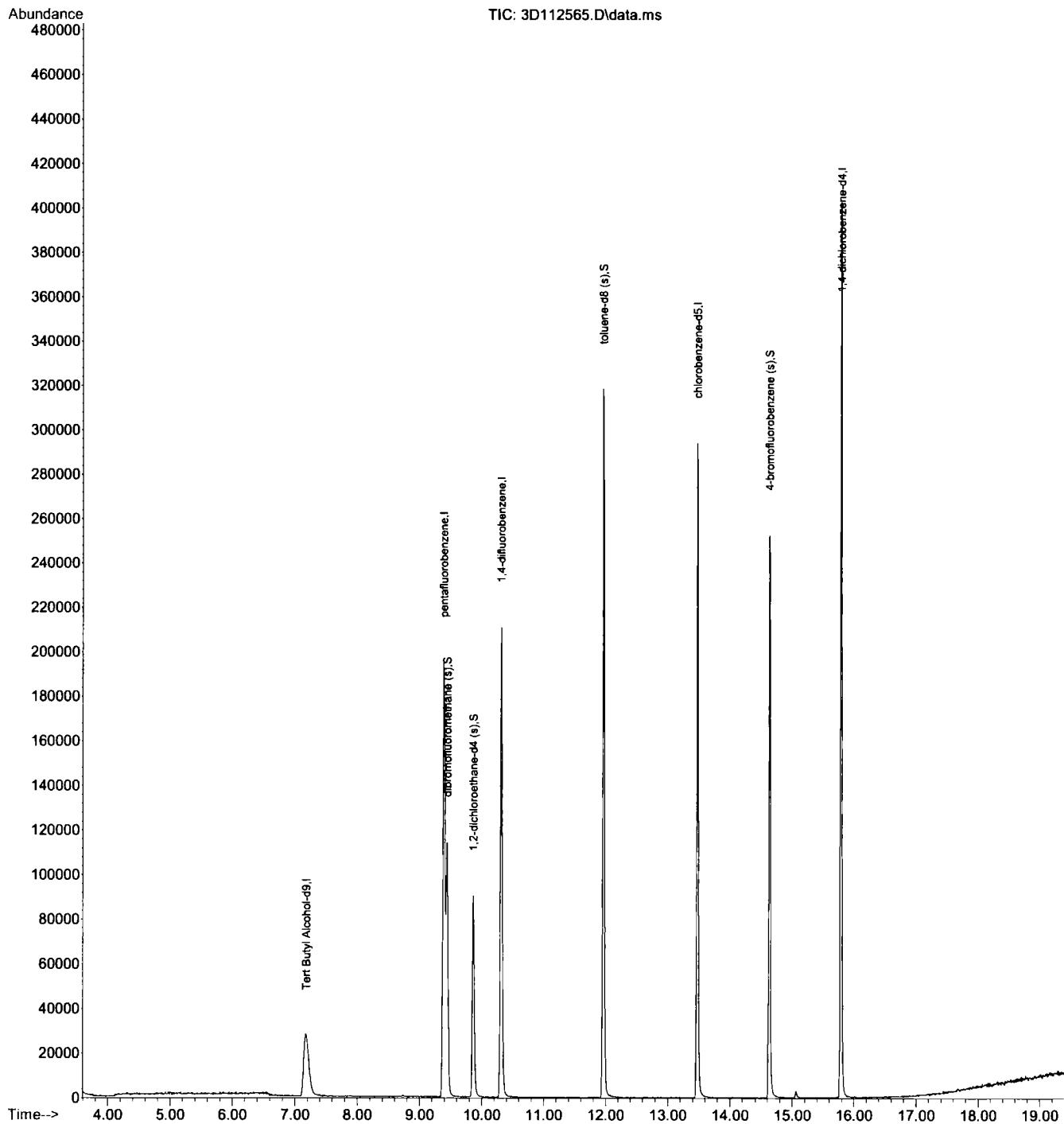
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	78922	56.34	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	112.68%
51) 1,2-dichloroethane-d4 (s)	9.863	65	72682	56.01	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	112.02%
80) toluene-d8 (s)	11.960	98	234935	51.09	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	102.18%
106) 4-bromofluorobenzene (s)	14.629	95	80531	47.66	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	95.32%

Target Compounds	Qvalue
(#)= qualifier out of range (m)= manual integration (+)= signals summed	

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112565.D
 Acq On : 12 Aug 2015 10:15 am
 Operator : ximenac
 Sample : mb2
 Misc : MS89457,V3D4823,5,,,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Aug 12 14:29:31 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112520.D
 Acq On : 11 Aug 2015 1:12 pm
 Operator : ximenac
 Sample : bs
 Misc : MS88280,V3D4822,5,,,1
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 11 16:10:50 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.178	65	107290	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	202270	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	233933	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	196021	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	124533	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
50) dibromofluoromethane (s)	9.438	113	85099	54.91	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	= 109.82%		
51) 1,2-dichloroethane-d4 (s)	9.858	65	77743	54.15	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	= 108.30%		
80) toluene-d8 (s)	11.960	98	263114	51.36	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	= 102.72%		
106) 4-bromofluorobenzene (s)	14.629	95	94978	49.57	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	= 99.14%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-dioxane	11.032	88	26388	1522.14	ug/L	96
3) tertiary butyl alcohol	7.278	59	57465	248.00	ug/L	87
8) chlorodifluoromethane	3.860	51	119080	43.37	ug/L	98
9) dichlorodifluoromethane	3.844	85	155036	42.17	ug/L	99
10) chloromethane	4.164	50	136902	46.83	ug/L	98
11) vinyl chloride	4.415	62	157814	53.83	ug/L	99
12) bromomethane	5.050	94	97837	51.65	ug/L	99
13) chloroethane	5.212	64	73255	58.99	ug/L	100
16) trichlorofluoromethane	5.679	101	167127	54.62	ug/L	99
18) ethyl ether	6.067	74	40832	54.03	ug/L	96
21) acrolein	6.319	56	161558	532.55	ug/L	99
22) 1,1-dichloroethene	6.486	61	134027	55.71	ug/L	96
23) acetone	6.533	43	25538	56.35	ug/L	90
24) allyl chloride	7.011	76	49193	65.83	ug/L #	85
25) acetonitrile	6.969	40	85772	569.14	ug/L	93
27) iodomethane	6.759	142	202177	59.10	ug/L	97
28) iso-butyl alcohol	9.679	41	28790	571.72	ug/L	94
29) carbon disulfide	6.890	76	362498	59.58	ug/L	99
30) methylene chloride	7.199	84	102677	56.95	ug/L	99
31) methyl acetate	6.995	43	54210	53.84	ug/L	98
32) methyl tert butyl ether	7.524	73	565686	105.92	ug/L	94
33) trans-1,2-dichloroethene	7.572	61	122774	55.93	ug/L	98
34) di-isopropyl ether	8.117	45	277261	49.11	ug/L	93
35) ethyl tert-butyl ether	8.584	59	279165	52.48	ug/L	98
36) 2-butanone	8.846	72	8206	56.89	ug/L	92
37) 1,1-dichloroethane	8.138	63	155048	59.45	ug/L	98
38) chloroprene	8.243	53	95155	49.85	ug/L	97
39) acrylonitrile	7.524	53	160542	307.57	ug/L	97
40) vinyl acetate	8.127	86	11654	60.15	ug/L	77
41) ethyl acetate	8.861	45	9049	45.61	ug/L #	20
42) 2,2-dichloropropane	8.877	77	155410	56.94	ug/L	99
43) cis-1,2-dichloroethene	8.872	96	92333	52.38	ug/L	97
44) propionitrile	8.935	54	107314	551.71	ug/L	95
45) methyl acrylate	8.940	55	75433	57.53	ug/L	99
46) bromochloromethane	9.186	128	48665	57.37	ug/L	95
47) tetrahydrofuran	9.228	42	25558	50.02	ug/L	99
48) chloroform	9.239	83	143053	56.89	ug/L	100
49) Tert-Butyl Formate	9.276	59	79464	54.28	ug/L #	96
52) freon 113	6.455	151	80877	54.60	ug/L	99
53) methacrylonitrile	9.118	41	39382	51.82	ug/L	97
54) 1,1,1-trichloroethane	9.496	97	156762	60.38	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112520.D
 Acq On : 11 Aug 2015 1:12 pm
 Operator : ximenac
 Sample : bs
 Misc : MS88280,V3D4822,5,,,1
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 11 16:10:50 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
57) tert-amyl methyl ether	9.968	73	276395	51.99	ug/L	97
59) epichlorohydrin	11.572	57	35410	275.73	ug/L	97
60) n-butyl alcohol	10.440	56	110794	2900.73	ug/L	99
61) cyclohexane	9.569	84	137794	50.30	ug/L	99
62) carbon tetrachloride	9.700	117	143116	63.26	ug/L	99
63) 1,1-dichloropropene	9.674	75	98363	58.38	ug/L	99
64) hexane	7.865	57	68294	45.38	ug/L	99
65) benzene	9.936	78	300665	54.70	ug/L	99
66) heptane	10.109	57	39107	45.74	ug/L	96
67) isopropyl acetate	9.858	43	143187	59.88	ug/L	96
68) 1,2-dichloroethane	9.952	62	91902	55.98	ug/L	98
69) trichloroethene	10.655	95	73027	54.95	ug/L	98
72) 2-nitropropane	11.436	41	26473	54.49	ug/L	91
73) 2-chloroethyl vinyl ether	11.446	63	253359	300.29	ug/L	98
74) methyl methacrylate	10.922	41	107868	57.40	ug/L	96
75) 1,2-dichloropropane	10.917	63	78705	54.78	ug/L	92
76) methylcyclohexane	10.875	83	130622	52.54	ug/L	96
77) dibromomethane	11.079	93	49016	57.99	ug/L	92
78) bromodichloromethane	11.210	83	98620	56.41	ug/L	98
79) cis-1,3-dichloropropene	11.666	75	115152	56.10	ug/L	96
81) 4-methyl-2-pentanone	11.766	58	29415	56.24	ug/L	99
82) toluene	12.033	92	171857	56.89	ug/L	100
83) 3-methyl-1-butanol	11.787	70	44704	1131.08	ug/L	97
84) trans-1,3-dichloropropene	12.233	75	100738	55.23	ug/L	96
85) ethyl methacrylate	12.227	69	79632	49.40	ug/L	97
86) 1,1,2-trichloroethane	12.442	83	56371	55.55	ug/L	96
87) 2-hexanone	12.621	58	22188	51.90	ug/L	89
89) tetrachloroethene	12.621	166	84824	50.82	ug/L	96
90) 1,3-dichloropropane	12.631	76	101486	53.36	ug/L	97
91) butyl acetate	12.694	56	42598	54.12	ug/L	90
92) 3,3-Dimethyl-1-Butanol	12.788	57	83134	441.89	ug/L	98
93) dibromochloromethane	12.893	129	86173	58.10	ug/L	99
94) 1,2-dibromoethane	13.045	107	68520	54.79	ug/L	97
95) n-Butyl Ether	13.423	57	283536	47.82	ug/L	98
96) chlorobenzene	13.507	112	195625	55.65	ug/L	100
97) 1,1,1,2-tetrachloroethane	13.564	131	90129	57.81	ug/L	98
98) ethylbenzene	13.559	91	314792	52.28	ug/L	98
99) m,p-xylene	13.664	106	250310	111.17	ug/L	98
100) o-xylene	14.083	106	130774	56.86	ug/L	97
101) styrene	14.094	104	203516	52.59	ug/L	99
103) bromoform	14.356	173	66275	57.27	ug/L	99
105) isopropylbenzene	14.424	105	344313	55.10	ug/L	99
107) bromobenzene	14.823	156	98388	53.34	ug/L	98
108) cyclohexanone	14.587	55	53873	284.81	ug/L	97
109) 1,1,2,2-tetrachloroethane	14.728	83	97454	53.89	ug/L	98
110) trans-1,4-dichloro-2-b...	14.770	53	18920	47.27	ug/L	98
111) 1,2,3-trichloropropene	14.802	110	23435	56.74	ug/L	97
112) n-propylbenzene	14.833	91	398245	58.53	ug/L	99
114) 2-chlorotoluene	14.980	126	88155	55.47	ug/L	98
115) 4-chlorotoluene	15.080	126	84081	54.48	ug/L	96
116) 1,3,5-trimethylbenzene	14.985	105	297645	55.97	ug/L	98
117) tert-butylbenzene	15.331	119	243005	55.45	ug/L	99
118) pentachloroethane	15.410	167	76314	63.86	ug/L	97
119) 1,2,4-trimethylbenzene	15.378	105	308226	57.73	ug/L	99
120) sec-butylbenzene	15.546	105	402481	57.48	ug/L	99
121) 1,3-dichlorobenzene	15.735	146	188957	55.10	ug/L	99
122) p-isopropyltoluene	15.662	119	349193	57.11	ug/L	99
123) 1,4-dichlorobenzene	15.814	146	197048	53.56	ug/L	100

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112520.D
 Acq On : 11 Aug 2015 1:12 pm
 Operator : ximenac
 Sample : bs
 Misc : MS88280,V3D4822,5,,,1
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 11 16:10:50 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

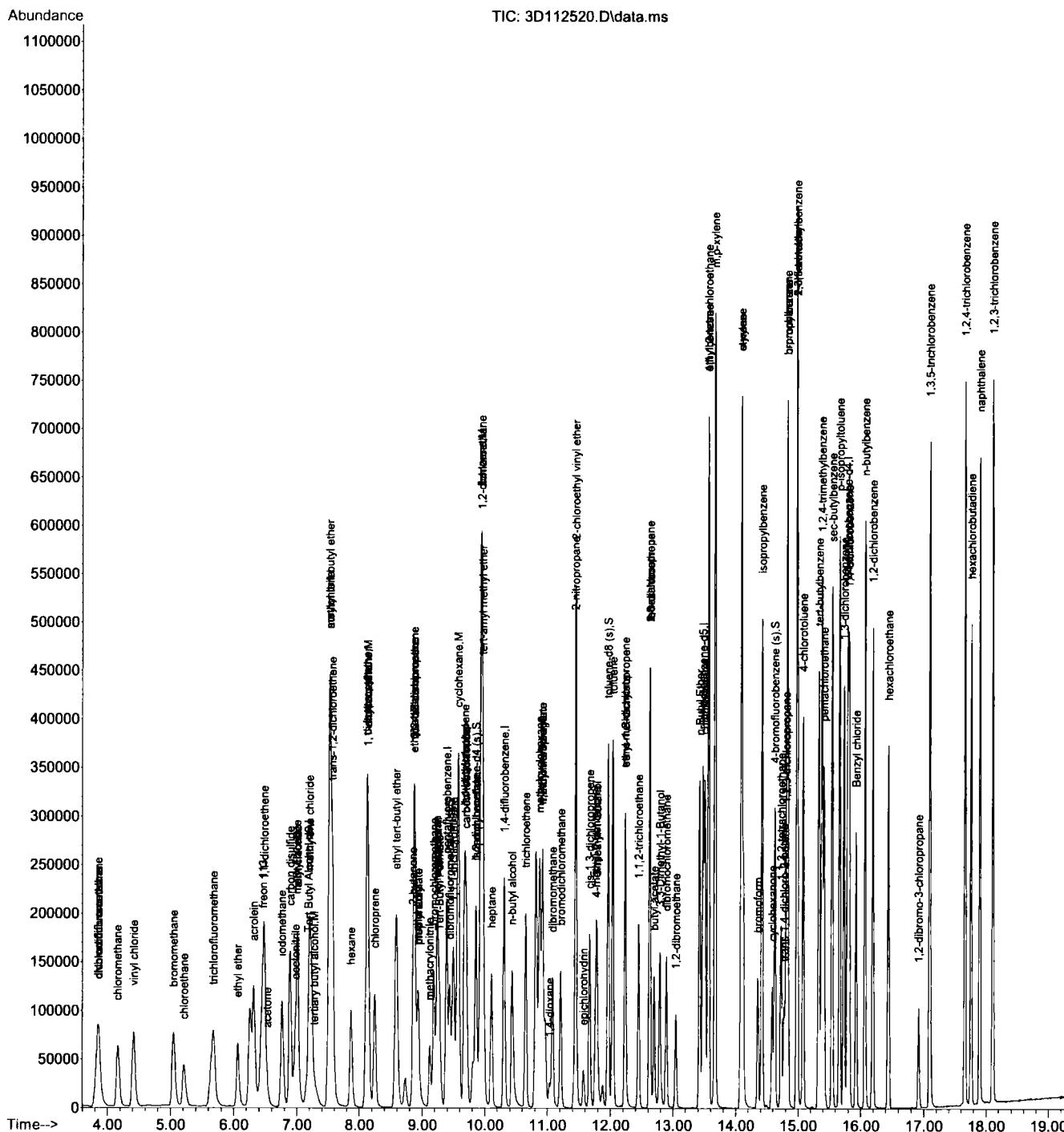
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.191	146	205920	55.70	ug/L	97
126) n-butylbenzene	16.065	92	181028	56.58	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.920	157	28959	53.60	ug/L	95
129) 1,3,5-trichlorobenzene	17.093	180	229423	58.44	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	230272	57.07	ug/L	96
131) hexachlorobutadiene	17.759	225	108359	54.25	ug/L	99
132) naphthalene	17.900	128	483762	56.60	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	240090	57.31	ug/L	98
134) hexachloroethane	16.443	119	72530	57.43	ug/L	98
135) Benzyl chloride	15.924	91	201481	56.09	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
Data File : 3D112520.D
Acq On : 11 Aug 2015 1:12 pm
Operator : ximenac
Sample : bs
Misc : MS88280,V3D4822,5,,,1
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Aug 11 16:10:50 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



M3D4810.M Tue Aug 11 16:12:35 2015 3D

Page : 4

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112545.D
 Acq On : 12 Aug 2015 12:38 am
 Operator : ximenac
 Sample : bs
 Misc : MS89468,V3D4823,5,,,1
 ALS Vial : 38 Sample Multiplier: 1

Quant Time: Aug 12 09:41:05 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	93840	500.00	ug/L	0.00
4) pentafluorobenzene	9.396	168	196820	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	222078	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	188570	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	114311	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.449	113	81749	54.21	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	= 108.42%		
51) 1,2-dichloroethane-d4 (s)	9.868	65	73671	52.74	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	= 105.48%		
80) toluene-d8 (s)	11.965	98	253218	52.06	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	= 104.12%		
106) 4-bromofluorobenzene (s)	14.629	95	89276	50.76	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	= 101.52%		
Target Compounds						
				Qvalue		
2) 1,4-dioxane	11.042	88	22454	1480.86	ug/L	95
3) tertiary butyl alcohol	7.288	59	47905	236.38	ug/L	92
8) chlorodifluoromethane	3.844	51	111846	41.86	ug/L	95
9) dichlorodifluoromethane	3.849	85	208243	58.21	ug/L	100
10) chloromethane	4.164	50	150580	52.94	ug/L	98
11) vinyl chloride	4.415	62	171832	60.23	ug/L	100
12) bromomethane	5.050	94	103954	56.40	ug/L	97
13) chloroethane	5.212	64	75325	62.34	ug/L	97
16) trichlorofluoromethane	5.674	101	179133	60.17	ug/L	97
18) ethyl ether	6.067	74	36739	49.96	ug/L	95
21) acrolein	6.318	56	150454	509.68	ug/L	98
22) 1,1-dichloroethene	6.486	61	124138	53.03	ug/L	99
23) acetone	6.533	43	22174	50.28	ug/L	87
24) allyl chloride	7.011	76	42716	58.75	ug/L	99
25) acetonitrile	6.984	40	77590	529.11	ug/L	87
27) iodomethane	6.764	142	181805	54.62	ug/L	98
28) iso-butyl alcohol	9.690	41	24898	508.13	ug/L	91
29) carbon disulfide	6.890	76	327219	55.27	ug/L	99
30) methylene chloride	7.205	84	93869	53.51	ug/L	97
31) methyl acetate	7.000	43	47084	48.06	ug/L	99
32) methyl tert butyl ether	7.530	73	516780	99.44	ug/L	94
33) trans-1,2-dichloroethene	7.577	61	113434	53.11	ug/L	98
34) di-isopropyl ether	8.127	45	253704	46.18	ug/L	94
35) ethyl tert-butyl ether	8.594	59	255067	49.28	ug/L	98
36) 2-butanone	8.851	72	7376	52.56	ug/L	88
37) 1,1-dichloroethane	8.143	63	143347	56.49	ug/L	100
38) chloroprene	8.253	53	89192	48.02	ug/L	98
39) acrylonitrile	7.530	53	142877	281.30	ug/L	96
40) vinyl acetate	8.127	86	10693	56.71	ug/L	96
41) ethyl acetate	8.867	45	7991	41.39	ug/L #	38
42) 2,2-dichloropropane	8.882	77	132446	49.87	ug/L	98
43) cis-1,2-dichloroethene	8.882	96	83967	48.95	ug/L	99
44) propionitrile	8.945	54	95957	506.98	ug/L	96
45) methyl acrylate	8.945	55	65257	51.15	ug/L	99
46) bromochloromethane	9.192	128	44522	53.94	ug/L	95
47) tetrahydrofuran	9.234	42	22560	45.38	ug/L	99
48) chloroform	9.244	83	131556	53.76	ug/L	98
49) Tert-Butyl Formate	9.276	59	70426	49.44	ug/L #	96
52) freon 113	6.450	151	81194	56.33	ug/L	99
53) methacrylonitrile	9.129	41	35332	47.78	ug/L	99

M3D4810.M Wed Aug 12 09:47:55 2015 T

Page: 1

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112545.D
 Acq On : 12 Aug 2015 12:38 am
 Operator : ximenac
 Sample : bs
 Misc : MS89468,V3D4823,5,,,1
 ALS Vial : 38 Sample Multiplier: 1

Quant Time: Aug 12 09:41:05 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
54) 1,1,1-trichloroethane	9.501	97	146818	58.11	ug/L	99
57) tert-amyl methyl ether	9.973	73	249752	48.28	ug/L	97
59) epichlorohydrin	11.572	57	30078	246.71	ug/L	98
60) n-butyl alcohol	10.445	56	94866	2616.30	ug/L	99
61) cyclohexane	9.574	84	130659	50.24	ug/L	100
62) carbon tetrachloride	9.705	117	135652	63.16	ug/L	99
63) 1,1-dichloropropene	9.679	75	91530	57.22	ug/L	97
64) hexane	7.870	57	62156	43.51	ug/L	97
65) benzene	9.941	78	279967	53.65	ug/L	99
66) heptane	10.109	57	36081	44.46	ug/L	96
67) isopropyl acetate	9.863	43	127029	55.96	ug/L	97
68) 1,2-dichloroethane	9.957	62	86716	55.64	ug/L	99
69) trichloroethene	10.660	95	67073	53.16	ug/L	98
72) 2-nitropropane	11.430	41	23895	51.80	ug/L	96
73) 2-chloroethyl vinyl ether	11.451	63	231880	289.50	ug/L	98
74) methyl methacrylate	10.927	41	94285	52.85	ug/L	95
75) 1,2-dichloropropane	10.922	63	74833	54.86	ug/L	96
76) methylcyclohexane	10.880	83	127691	54.10	ug/L	94
77) dibromomethane	11.084	93	45381	56.55	ug/L	89
78) bromodichloromethane	11.215	83	91949	55.40	ug/L	100
79) cis-1,3-dichloropropene	11.677	75	105078	53.93	ug/L	95
81) 4-methyl-2-pentanone	11.766	58	26773	53.92	ug/L	92
82) toluene	12.039	92	159793	55.72	ug/L	98
83) 3-methyl-1-butanol	11.787	70	37380	996.26	ug/L	96
84) trans-1,3-dichloropropene	12.238	75	91249	52.70	ug/L	94
85) ethyl methacrylate	12.227	69	72242	47.20	ug/L	98
86) 1,1,2-trichloroethane	12.448	83	52270	54.26	ug/L	96
87) 2-hexanone	12.621	58	18597	45.82	ug/L	97
89) tetrachloroethene	12.626	166	84616	52.70	ug/L	99
90) 1,3-dichloropropane	12.631	76	93663	51.19	ug/L	95
91) butyl acetate	12.694	56	36841	48.65	ug/L	94
92) 3,3-Dimethyl-1-Butanol	12.794	57	67406	372.44	ug/L	96
93) dibromochloromethane	12.893	129	78537	55.05	ug/L	99
94) 1,2-dibromoethane	13.045	107	62690	52.10	ug/L	96
95) n-Butyl Ether	13.423	57	254755	44.66	ug/L	98
96) chlorobenzene	13.507	112	180854	53.48	ug/L	99
97) 1,1,1,2-tetrachloroethane	13.564	131	83757	55.85	ug/L	99
98) ethylbenzene	13.559	91	293168	50.61	ug/L	99
99) m,p-xylene	13.669	106	231968	107.09	ug/L	99
100) o-xylene	14.083	106	120260	54.35	ug/L	98
101) styrene	14.094	104	184337	49.52	ug/L	97
103) bromoform	14.361	173	59621	53.56	ug/L	99
105) isopropylbenzene	14.424	105	318960	55.60	ug/L	99
107) bromobenzene	14.823	156	89959	53.13	ug/L	98
108) cyclohexanone	14.587	55	67342	367.69	ug/L	99
109) 1,1,2,2-tetrachloroethane	14.728	83	87461	52.68	ug/L	99
110) trans-1,4-dichloro-2-b...	14.770	53	14374	39.12	ug/L	98
111) 1,2,3-trichloropropane	14.802	110	20389	53.78	ug/L	92
112) n-propylbenzene	14.833	91	363649	58.22	ug/L	100
114) 2-chlorotoluene	14.980	126	79954	54.80	ug/L	97
115) 4-chlorotoluene	15.080	126	76073	53.70	ug/L	96
116) 1,3,5-trimethylbenzene	14.985	105	275161	56.37	ug/L	100
117) tert-butylbenzene	15.331	119	226383	56.28	ug/L	100
118) pentachloroethane	15.410	167	66238	60.38	ug/L	97
119) 1,2,4-trimethylbenzene	15.378	105	280691	57.27	ug/L	98
120) sec-butylbenzene	15.546	105	372952	58.02	ug/L	100
121) 1,3-dichlorobenzene	15.735	146	172633	54.84	ug/L	98
122) p-isopropyltoluene	15.662	119	316285	56.35	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112545.D
 Acq On : 12 Aug 2015 12:38 am
 Operator : ximenac
 Sample : bs
 Misc : MS89468,V3D4823,5,,,1
 ALS Vial : 38 Sample Multiplier: 1

Quant Time: Aug 12 09:41:05 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
123) 1,4-dichlorobenzene	15.814	146	176213	52.18	ug/L	98
124) 1,2-dichlorobenzene	16.191	146	186915	55.08	ug/L	99
126) n-butylbenzene	16.065	92	159806	54.42	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.920	157	24277	48.95	ug/L	97
129) 1,3,5-trichlorobenzene	17.093	180	201841	56.02	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	205029	55.35	ug/L	97
131) hexachlorobutadiene	17.759	225	98241	53.58	ug/L	98
132) naphthalene	17.900	128	423480	53.97	ug/L	99
133) 1,2,3-trichlorobenzene	18.110	180	215311	55.99	ug/L	98
134) hexachloroethane	16.443	119	66356	57.24	ug/L	98
135) Benzyl chloride	15.924	91	146174	44.33	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

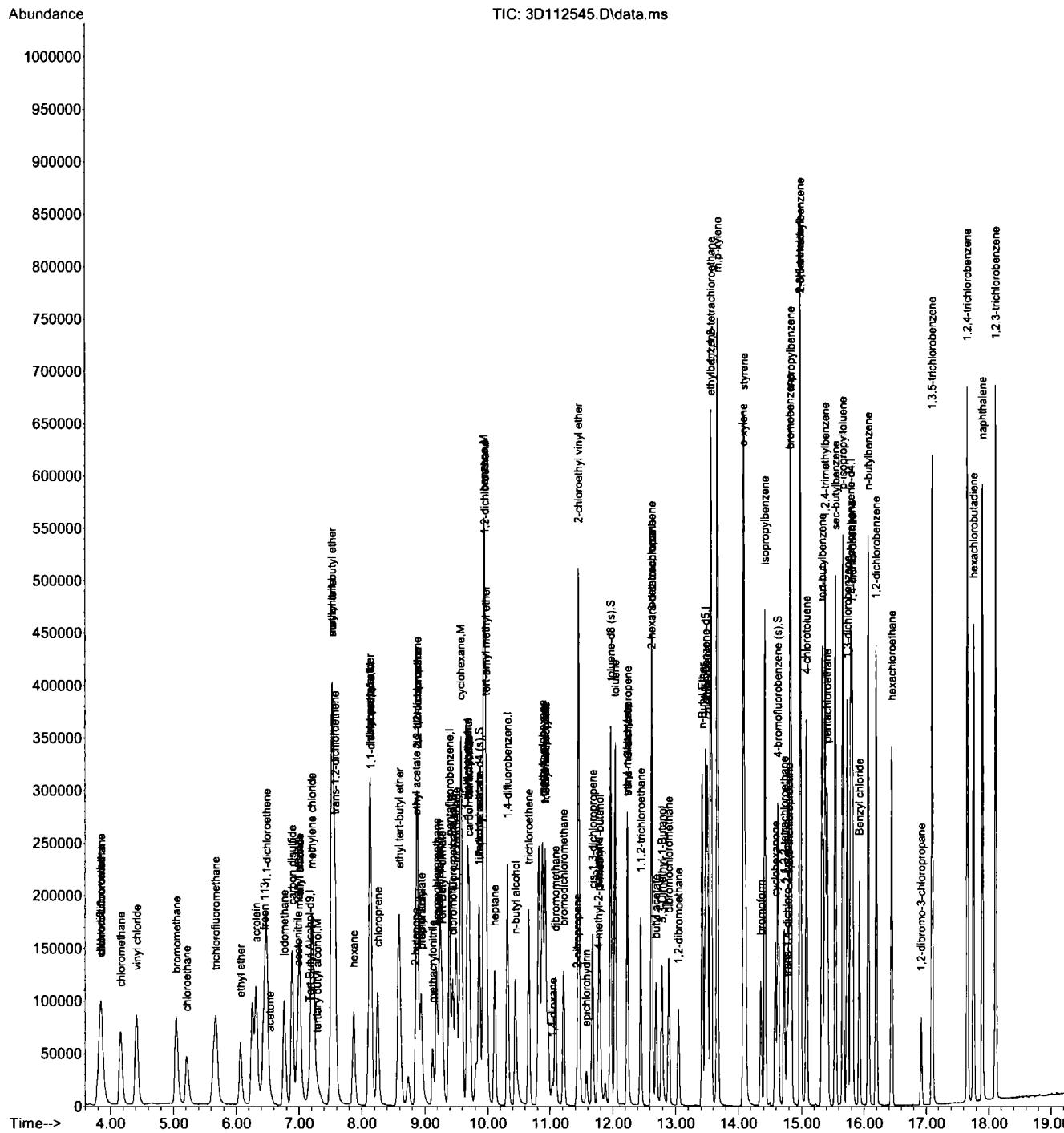
7.3.2

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Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112545.D
 Acq On : 12 Aug 2015 12:38 am
 Operator : ximenac
 Sample : bs
 Misc : MS89468, V3D4823, 5, , , 1
 ALS Vial : 38 Sample Multiplier: 1

Quant Time: Aug 12 09:41:05 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112570.D
 Acq On : 12 Aug 2015 12:33 pm
 Operator : ximenac
 Sample : bs
 Misc : MS89457,V3D4824,5,,,1
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Aug 12 14:31:21 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.168	65	97953	500.00	ug/L	-0.01
4) pentafluorobenzene	9.391	168	207104	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	238560	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	200201	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	122169	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.438	113	87153	54.92	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 109.84%	
51) 1,2-dichloroethane-d4 (s)	9.858	65	78891	53.67	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 107.34%	
80) toluene-d8 (s)	11.960	98	268063	51.31	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 102.62%	
106) 4-bromofluorobenzene (s)	14.629	95	94955	50.52	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 101.04%	
Target Compounds						
2) 1,4-dioxane	11.037	88	23107	1459.94	ug/L	99
3) tertiary butyl alcohol	7.289	59	49500	233.99	ug/L	93
8) chlorodifluoromethane	3.849	51	153661	54.65	ug/L	98
9) dichlorodifluoromethane	3.839	85	193726	51.47	ug/L	99
10) chloromethane	4.158	50	142073	47.47	ug/L	99
11) vinyl chloride	4.415	62	165704	55.20	ug/L	99
12) bromomethane	5.050	94	100041	51.58	ug/L	98
13) chloroethane	5.218	64	73785	58.03	ug/L	96
16) trichlorofluoromethane	5.674	101	176168	56.24	ug/L	99
18) ethyl ether	6.067	74	39493	51.04	ug/L	94
21) acrolein	6.319	56	140427	452.09	ug/L	100
22) 1,1-dichloroethene	6.481	61	129346	52.51	ug/L	96
23) acetone	6.528	43	20998	45.25	ug/L	85
24) allyl chloride	7.011	76	45363	59.29	ug/L	91
25) acetonitrile	6.969	40	80633	522.56	ug/L	88
27) iodomethane	6.764	142	197384	56.35	ug/L	99
28) iso-butyl alcohol	9.685	41	25332	491.31	ug/L	96
29) carbon disulfide	6.890	76	354175	56.85	ug/L	99
30) methylene chloride	7.205	84	98215	53.21	ug/L	97
31) methyl acetate	6.995	43	51449	49.91	ug/L	97
32) methyl tert butyl ether	7.530	73	272061	49.75	ug/L	87
33) trans-1,2-dichloroethene	7.572	61	117711	52.37	ug/L	97
34) di-isopropyl ether	8.117	45	277372	47.98	ug/L	99
35) ethyl tert-butyl ether	8.584	59	281106	51.61	ug/L	97
36) 2-butanone	8.851	72	7348	49.76	ug/L	92
37) 1,1-dichloroethane	8.138	63	147764	55.34	ug/L	99
38) chloroprene	8.248	53	98805	50.55	ug/L	97
39) acrylonitrile	7.524	53	141114	264.04	ug/L	98
40) vinyl acetate	8.127	86	11261	56.76	ug/L	96
41) ethyl acetate	8.856	45	8631	42.48	ug/L	91
42) 2,2-dichloropropane	8.877	77	153682	54.99	ug/L	97
43) cis-1,2-dichloroethene	8.877	96	88672	49.13	ug/L	100
44) propionitrile	8.935	54	96747	485.78	ug/L	95
45) methyl acrylate	8.940	55	66725	49.70	ug/L	98
46) bromochloromethane	9.186	128	46574	53.62	ug/L	96
47) tetrahydrofuran	9.223	42	22664	43.32	ug/L	99
48) chloroform	9.244	83	138603	53.83	ug/L	98
49) Tert-Butyl Formate	9.276	59	73660	49.14	ug/L #	97
52) freon 113	6.460	151	87838	57.91	ug/L	97
53) methacrylonitrile	9.124	41	35867	46.10	ug/L	97
54) 1,1,l-trichloroethane	9.496	97	149220	56.13	ug/L	96

M3D4810.M Wed Aug 12 14:32:49 2015 3D

Page: 1

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112570.D
 Acq On : 12 Aug 2015 12:33 pm
 Operator : ximenac
 Sample : bs
 Misc : MS89457,V3D4824,5,,,1
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Aug 12 14:31:21 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
57) tert-amyl methyl ether	9.973	73	269551	49.52	ug/L	97
59) epichlorohydrin	11.572	57	29377	224.31	ug/L	99
60) n-butyl alcohol	10.440	56	93101	2390.23	ug/L	97
61) cyclohexane	9.569	84	141274	50.57	ug/L	95
62) carbon tetrachloride	9.700	117	138744	60.13	ug/L	99
63) 1,1-dichloropropene	9.674	75	96108	55.93	ug/L	99
64) hexane	7.870	57	153694	100.15	ug/L	99
65) benzene	9.936	78	289842	51.71	ug/L	98
66) heptane	10.109	57	48940	56.14	ug/L	98
67) isopropyl acetate	9.858	43	114685	47.03	ug/L	98
68) 1,2-dichloroethane	9.952	62	89748	53.61	ug/L	99
69) trichloroethene	10.655	95	70352	51.91	ug/L	98
72) 2-nitropropane	11.431	41	23966	48.37	ug/L	100
73) 2-chloroethyl vinyl ether	11.446	63	214994	249.88	ug/L	99
74) methyl methacrylate	10.922	41	98131	51.20	ug/L	95
75) 1,2-dichloropropane	10.917	63	76357	52.11	ug/L	95
76) methylcyclohexane	10.875	83	148370	58.52	ug/L	96
77) dibromomethane	11.079	93	47408	55.00	ug/L	95
78) bromodichloromethane	11.210	83	95889	53.78	ug/L	97
79) cis-1,3-dichloropropene	11.672	75	110107	52.61	ug/L	98
81) 4-methyl-2-pentanone	11.761	58	26137	49.00	ug/L	96
82) toluene	12.033	92	166153	53.94	ug/L	99
83) 3-methyl-1-butanol	11.787	70	36795	912.92	ug/L	99
84) trans-1,3-dichloropropene	12.233	75	96369	51.81	ug/L	98
85) ethyl methacrylate	12.227	69	72670	44.20	ug/L	99
86) 1,1,2-trichloroethane	12.442	83	53697	51.89	ug/L	97
87) 2-hexanone	12.621	58	18504	42.44	ug/L	98
89) tetrachloroethene	12.626	166	82940	48.66	ug/L	100
90) 1,3-dichloropropane	12.626	76	95533	49.18	ug/L	96
91) butyl acetate	12.694	56	37223	46.30	ug/L	93
92) 3,3-Dimethyl-1-Butanol	12.788	57	73709	383.61	ug/L	97
93) dibromochloromethane	12.893	129	81309	53.68	ug/L	98
94) 1,2-dibromoethane	13.045	107	64561	50.54	ug/L	99
95) n-Butyl Ether	13.423	57	264688	43.71	ug/L	98
96) chlorobenzene	13.507	112	188969	52.63	ug/L	100
97) 1,1,1,2-tetrachloroethane	13.564	131	86382	54.25	ug/L	99
98) ethylbenzene	13.559	91	303312	49.32	ug/L	97
99) m,p-xylene	13.664	106	239234	104.03	ug/L	98
100) o-xylene	14.083	106	124621	53.05	ug/L	96
101) styrene	14.094	104	191744	48.52	ug/L	98
103) bromoform	14.356	173	60581	51.26	ug/L	99
105) isopropylbenzene	14.424	105	327663	53.45	ug/L	99
107) bromobenzene	14.823	156	93520	51.68	ug/L	98
108) cyclohexanone	14.587	55	26832	172.04	ug/L	95
109) 1,1,2,2-tetrachloroethane	14.728	83	88125	49.67	ug/L	100
110) trans-1,4-dichloro-2-b...	14.770	53	16262	41.41	ug/L	96
111) 1,2,3-trichloropropane	14.802	110	20989	51.80	ug/L	93
112) n-propylbenzene	14.833	91	377762	56.59	ug/L	100
114) 2-chlorotoluene	14.980	126	83572	53.60	ug/L	97
115) 4-chlorotoluene	15.080	126	80845	53.40	ug/L	95
116) 1,3,5-trimethylbenzene	14.985	105	287236	55.06	ug/L	97
117) tert-butylbenzene	15.331	119	236188	54.94	ug/L	98
118) pentachloroethane	15.410	167	74299	63.37	ug/L	96
119) 1,2,4-trimethylbenzene	15.379	105	290264	55.42	ug/L	98
120) sec-butylbenzene	15.546	105	392113	57.08	ug/L	99
121) 1,3-dichlorobenzene	15.730	146	180340	53.60	ug/L	99
122) p-isopropyltoluene	15.662	119	335368	55.91	ug/L	99
123) 1,4-dichlorobenzene	15.814	146	186307	51.62	ug/L	100

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112570.D
 Acq On : 12 Aug 2015 12:33 pm
 Operator : ximenac
 Sample : bs
 Misc : MS89457,V3D4824,5,,,1
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Aug 12 14:31:21 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

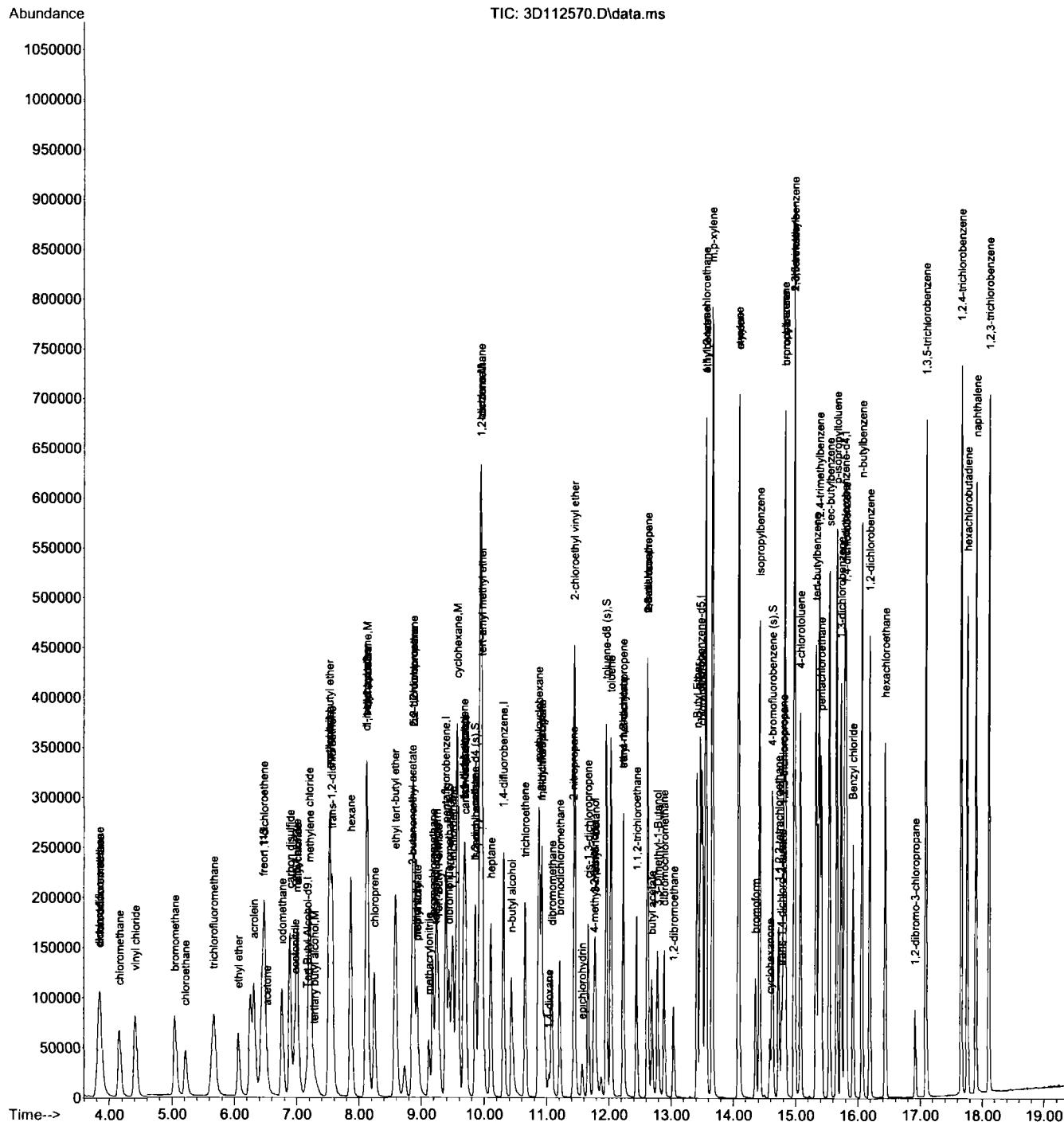
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.191	146	195325	53.85	ug/L	98
126) n-butylbenzene	16.065	92	173115	55.16	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.925	157	25560	48.22	ug/L	95
129) 1,3,5-trichlorobenzene	17.093	180	220676	57.30	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	225321	56.92	ug/L	98
131) hexachlorobutadiene	17.759	225	108818	55.54	ug/L	97
132) naphthalene	17.900	128	443635	52.91	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	229725	55.90	ug/L	99
134) hexachloroethane	16.438	119	70374	56.80	ug/L	98
135) Benzyl chloride	15.924	91	178287	50.59	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
Data File : 3D112570.D
Acq On : 12 Aug 2015 12:33 pm
Operator : ximenac
Sample : bs
Misc : MS89457,V3D4824,5,,,1
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Aug 12 14:31:21 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112526.D
 Acq On : 11 Aug 2015 4:03 pm
 Operator : ximenac
 Sample : jc1106-13ms
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 12 09:27:51 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.168	65	109061	500.00	ug/L	-0.01
4) pentafluorobenzene	9.391	168	200873	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	227802	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	191425	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	118628	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	84366	54.82	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	= 109.64%		
51) 1,2-dichloroethane-d4 (s)	9.863	65	75945	53.27	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	= 106.54%		
80) toluene-d8 (s)	11.960	98	257948	51.70	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	= 103.40%		
106) 4-bromofluorobenzene (s)	14.629	95	91571	50.17	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	= 100.34%		
Target Compounds						
				Qvalue		
2) 1,4-dioxane	11.037	88	28995	1645.36	ug/L	97
3) tertiary butyl alcohol	7.288	59	62710	266.24	ug/L	94
8) chlorodifluoromethane	3.849	51	131260	48.13	ug/L	97
9) dichlorodifluoromethane	3.849	85	131473	36.01	ug/L	100
10) chloromethane	4.158	50	115576	39.81	ug/L	98
11) vinyl chloride	4.415	62	126943	43.60	ug/L	99
12) bromomethane	5.050	94	84879	45.12	ug/L	98
13) chloroethane	5.212	64	55676	45.15	ug/L	95
16) trichlorofluoromethane	5.679	101	145920	48.03	ug/L	98
18) ethyl ether	6.062	74	42546	56.69	ug/L	97
21) acrolein	6.319	56	155848	517.30	ug/L	99
22) 1,1-dichloroethene	6.486	61	129425	54.17	ug/L	96
23) acetone	6.539	43	26731	59.39	ug/L	92
24) allyl chloride	7.011	76	42264	56.95	ug/L	98
25) acetonitrile	6.979	40	91362	610.45	ug/L	91
27) iodomethane	6.764	142	207875	61.19	ug/L	99
28) iso-butyl alcohol	9.685	41	30786	615.61	ug/L	87
29) carbon disulfide	6.890	76	345032	57.10	ug/L	98
30) methylene chloride	7.199	84	106207	59.32	ug/L	99
31) methyl acetate	6.995	43	60840	60.85	ug/L	98
32) methyl tert butyl ether	7.530	73	304660	57.44	ug/L	98
33) trans-1,2-dichloroethene	7.572	61	124362	57.05	ug/L	100
34) di-isopropyl ether	8.122	45	286627	51.12	ug/L	91
35) ethyl tert-butyl ether	8.584	59	289842	54.87	ug/L	99
36) 2-butanone	8.846	72	8398	58.63	ug/L	93
37) 1,1-dichloroethane	8.138	63	157561	60.84	ug/L	100
38) chloroprene	8.243	53	89458	47.19	ug/L	96
39) acrylonitrile	7.524	53	167940	323.98	ug/L	98
40) vinyl acetate	8.122	86	11588	60.22	ug/L	87
41) ethyl acetate	8.861	45	9556	48.50	ug/L #	24
42) 2,2-dichloropropane	8.877	77	157191	57.99	ug/L	99
43) cis-1,2-dichloroethene	8.872	96	97111	55.47	ug/L	95
44) propionitrile	8.935	54	124100	642.45	ug/L	100
45) methyl acrylate	8.935	55	79232	60.85	ug/L	98
46) bromochloromethane	9.186	128	51313	60.91	ug/L	95
47) tetrahydrofuran	9.228	42	25820	50.88	ug/L	94
48) chloroform	9.244	83	144645	57.92	ug/L	98
49) Tert-Butyl Formate	9.270	59	69994	48.14	ug/L #	96
52) freon 113	6.444	151	74338	50.53	ug/L	98
53) methacrylonitrile	9.124	41	41340	54.78	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112526.D
 Acq On : 11 Aug 2015 4:03 pm
 Operator : ximenac
 Sample : jc1106-13ms
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 12 09:27:51 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
54) 1,1,1-trichloroethane	9.496	97	153086	59.37	ug/L	99
57) tert-amyl methyl ether	9.968	73	285704	54.11	ug/L	95
59) epichlorohydrin	11.572	57	33771	270.04	ug/L	99
60) n-butyl alcohol	10.434	56	122662	3297.88	ug/L	99
61) cyclohexane	9.569	84	133619	50.09	ug/L	97
62) carbon tetrachloride	9.700	117	131538	59.70	ug/L	98
63) 1,1-dichloropropene	9.669	75	89007	54.24	ug/L	99
64) hexane	7.865	57	62930	42.94	ug/L	97
65) benzene	9.936	78	298345	55.74	ug/L	98
66) heptane	10.104	57	40051	48.11	ug/L	98
67) isopropyl acetate	9.858	43	133566	57.36	ug/L	99
68) 1,2-dichloroethane	9.947	62	94610	59.18	ug/L	99
69) trichloroethene	10.655	95	71432	55.20	ug/L	95
72) 2-nitropropane	11.425	41	17785	37.59	ug/L	99
74) methyl methacrylate	10.917	41	104796	57.26	ug/L	99
75) 1,2-dichloropropane	10.917	63	79487	56.81	ug/L	94
76) methylcyclohexane	10.875	83	124108	51.26	ug/L	95
77) dibromomethane	11.079	93	51968	63.13	ug/L	93
78) bromodichloromethane	11.210	83	99138	58.23	ug/L	98
79) cis-1,3-dichloropropene	11.666	75	116188	58.13	ug/L	98
81) 4-methyl-2-pentanone	11.761	58	31981	62.79	ug/L	96
82) toluene	12.033	92	170956	58.12	ug/L	98
83) 3-methyl-1-butanol	11.782	70	51458	1337.01	ug/L	96
84) trans-1,3-dichloropropene	12.233	75	107380	60.45	ug/L	97
85) ethyl methacrylate	12.222	69	84472	53.81	ug/L	99
86) 1,1,2-trichloroethane	12.442	83	58312	59.01	ug/L	96
87) 2-hexanone	12.621	58	24591	59.06	ug/L	93
89) tetrachloroethene	12.621	166	80258	49.24	ug/L	98
90) 1,3-dichloropropane	12.626	76	104543	56.29	ug/L	96
91) butyl acetate	12.694	56	45521	59.22	ug/L	94
92) 3,3-Dimethyl-1-Butanol	12.788	57	114967	625.76	ug/L	99
93) dibromochloromethane	12.893	129	88282	60.95	ug/L	96
94) 1,2-dibromoethane	13.045	107	71420	58.47	ug/L	99
95) n-Butyl Ether	13.423	57	285198	49.25	ug/L	98
96) chlorobenzene	13.507	112	195854	57.05	ug/L	98
97) 1,1,1,2-tetrachloroethane	13.564	131	93304	61.29	ug/L	99
98) ethylbenzene	13.559	91	316316	53.80	ug/L	99
99) m,p-xylene	13.664	106	248269	112.91	ug/L	98
100) o-xylene	14.083	106	129181	57.51	ug/L	99
101) styrene	14.094	104	204004	53.99	ug/L	98
103) bromoform	14.356	173	67934	60.12	ug/L	99
105) isopropylbenzene	14.424	105	336418	56.51	ug/L	99
107) bromobenzene	14.823	156	100408	57.15	ug/L	97
108) cyclohexanone	14.587	55	31519	196.43	ug/L	97
109) 1,1,2,2-tetrachloroethane	14.728	83	104559	60.69	ug/L	99
110) trans-1,4-dichloro-2-b...	14.770	53	19881	52.14	ug/L	94
111) 1,2,3-trichloropropane	14.802	110	24032	61.08	ug/L	97
112) n-propylbenzene	14.833	91	364072	56.17	ug/L	100
114) 2-chlorotoluene	14.980	126	88376	58.37	ug/L	96
115) 4-chlorotoluene	15.080	126	84612	57.56	ug/L	95
116) 1,3,5-trimethylbenzene	14.985	105	298400	58.90	ug/L	98
117) tert-butylbenzene	15.331	119	234138	56.09	ug/L	98
118) pentachloroethane	15.410	167	78446	68.91	ug/L	98
119) 1,2,4-trimethylbenzene	15.379	105	294630	57.93	ug/L	99
120) sec-butylbenzene	15.546	105	387355	58.07	ug/L	99
121) 1,3-dichlorobenzene	15.730	146	190463	58.30	ug/L	99
122) p-isopropyltoluene	15.662	119	333891	57.33	ug/L	100
123) 1,4-dichlorobenzene	15.814	146	194167	55.40	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112526.D
 Acq On : 11 Aug 2015 4:03 pm
 Operator : ximenac
 Sample : jc1106-13ms
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 12 09:27:51 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

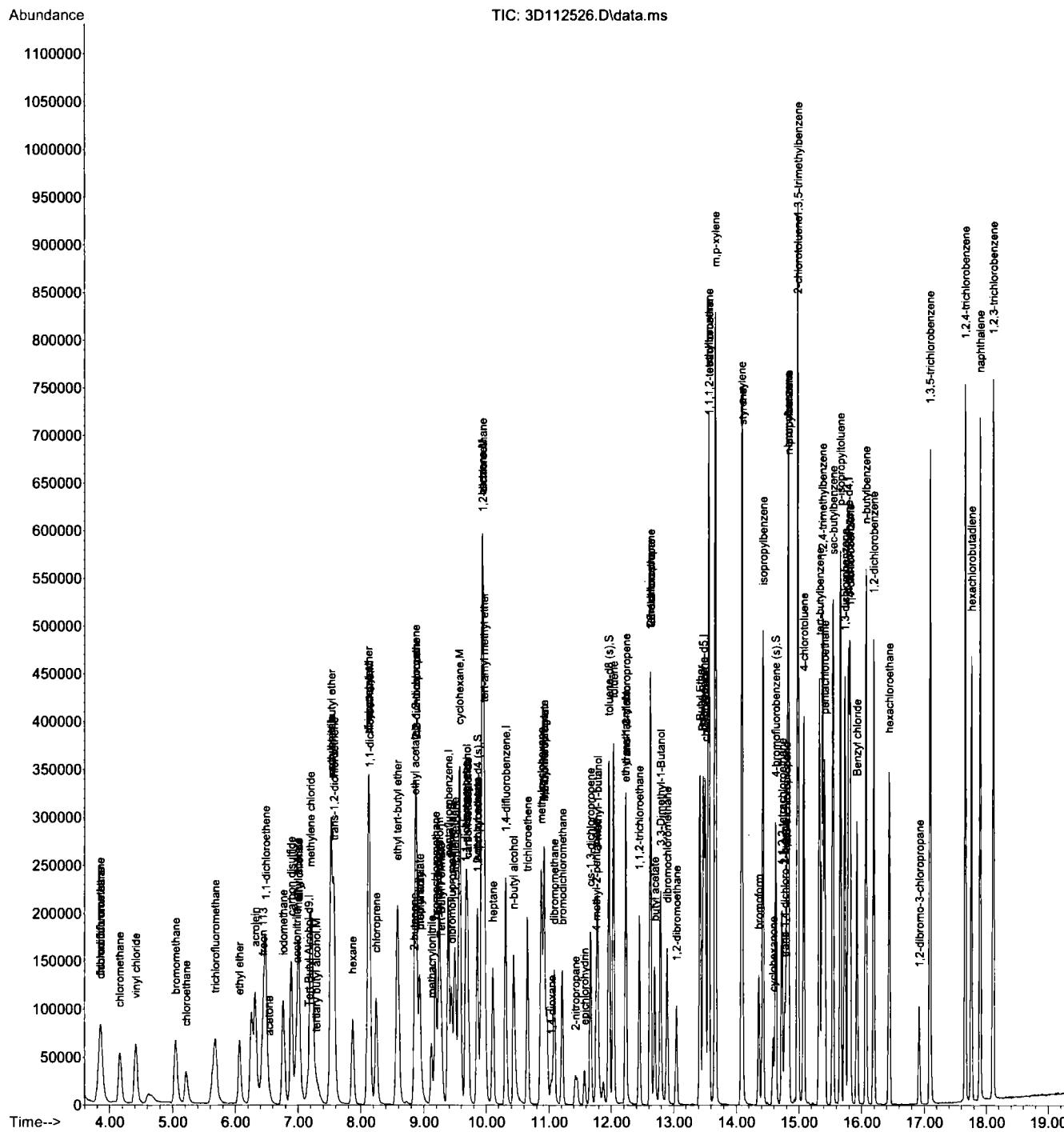
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.191	146	206200	58.55	ug/L	99
126) n-butylbenzene	16.065	92	169458	55.60	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.925	157	29809	57.91	ug/L	93
129) 1,3,5-trichlorobenzene	17.093	180	220521	58.97	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	230168	59.88	ug/L	97
131) hexachlorobutadiene	17.759	225	100067	52.59	ug/L	99
132) naphthalene	17.895	128	505438	62.08	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	242637	60.80	ug/L	98
134) hexachloroethane	16.438	119	70442	58.55	ug/L	98
135) Benzyl chloride	15.924	91	205038	59.92	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
Data File : 3D112526.D
Acq On : 11 Aug 2015 4:03 pm
Operator : ximenac
Sample : jc1106-13ms
Misc : MS89468,V3D4822,5,,,1
ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 12 09:27:51 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112527.D
 Acq On : 11 Aug 2015 4:30 pm
 Operator : ximenac
 Sample : jc1106-13msd
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 12 09:27:45 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.178	65	106157	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	209488	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	234394	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	194946	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	117427	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	86334	53.79	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	=	107.58%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	77471	52.10	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	=	104.20%	
80) toluene-d8 (s)	11.960	98	263504	51.33	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	=	102.66%	
106) 4-bromofluorobenzene (s)	14.629	95	92277	51.07	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	=	102.14%	
Target Compounds						
					Qvalue	
2) 1,4-dioxane	11.037	88	27251	1588.70	ug/L	92
3) tertiary butyl alcohol	7.294	59	62783	273.84	ug/L	95
8) chlorodifluoromethane	3.860	51	131503	46.24	ug/L	98
9) dichlorodifluoromethane	3.844	85	133435	35.04	ug/L	95
10) chloromethane	4.164	50	115335	38.10	ug/L	99
11) vinyl chloride	4.421	62	129232	42.56	ug/L	98
12) bromomethane	5.050	94	84829	43.24	ug/L	98
13) chloroethane	5.212	64	54043	42.02	ug/L	99
16) trichlorofluoromethane	5.679	101	144785	45.69	ug/L	97
18) ethyl ether	6.072	74	42484	54.28	ug/L	91
21) acrolein	6.319	56	155482	494.86	ug/L	99
22) 1,1-dichloroethene	6.486	61	125009	50.17	ug/L	95
23) acetone	6.533	43	24977	53.21	ug/L	84
24) allyl chloride	7.016	76	42273	54.62	ug/L	96
25) acetonitrile	6.974	40	90048	576.93	ug/L	91
27) iodomethane	6.764	142	206541	58.30	ug/L	99
28) iso-butyl alcohol	9.690	41	30622	587.15	ug/L	94
29) carbon disulfide	6.890	76	337722	53.60	ug/L	99
30) methylene chloride	7.199	84	105405	56.45	ug/L	99
31) methyl acetate	6.995	43	59945	57.49	ug/L	98
32) methyl tert butyl ether	7.535	73	304001	54.96	ug/L	99
33) trans-1,2-dichloroethene	7.577	61	122979	54.09	ug/L	99
34) di-isopropyl ether	8.122	45	287467	49.16	ug/L	96
35) ethyl tert-butyl ether	8.594	59	292254	53.05	ug/L	98
36) 2-butanone	8.851	72	8518	57.02	ug/L	86
37) 1,1-dichloroethane	8.143	63	155121	57.43	ug/L	99
38) chloroprene	8.248	53	90156	45.60	ug/L	97
39) acrylonitrile	7.530	53	168270	311.27	ug/L	95
40) vinyl acetate	8.127	86	11674	58.17	ug/L	79
41) ethyl acetate	8.861	45	9580	46.62	ug/L	62
42) 2,2-dichloropropane	8.888	77	154423	54.63	ug/L	97
43) cis-1,2-dichloroethene	8.877	96	95982	52.57	ug/L	99
44) propionitrile	8.935	54	123881	614.94	ug/L	95
45) methyl acrylate	8.945	55	78803	58.03	ug/L	98
46) bromochloromethane	9.186	128	52198	59.41	ug/L	97
47) tetrahydrofuran	9.234	42	25432	48.06	ug/L	96
48) chloroform	9.244	83	145612	55.91	ug/L	99
49) Tert-Butyl Formate	9.276	59	58356	38.49	ug/L #	97
52) freon 113	6.455	151	70632	46.04	ug/L	92
53) methacrylonitrile	9.124	41	40749	51.77	ug/L	98

M3D4810.M Wed Aug 12 09:47:46 2015 T

Page: 1

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112527.D
 Acq On : 11 Aug 2015 4:30 pm
 Operator : ximenac
 Sample : jc1106-13msd
 Misc : MS89468,V3D4822,5,,,1
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 12 09:27:45 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
54) 1,1,1-trichloroethane	9.501	97	150897	56.12	ug/L	99
57) tert-amyl methyl ether	9.973	73	285610	51.87	ug/L	97
59) epichlorohydrin	11.572	57	32481	252.42	ug/L	99
60) n-butyl alcohol	10.440	56	121033	3162.57	ug/L	99
61) cyclohexane	9.574	84	133372	48.59	ug/L	97
62) carbon tetrachloride	9.706	117	129099	56.95	ug/L	99
63) 1,1-dichloropropene	9.674	75	89346	52.92	ug/L	98
64) hexane	7.870	57	64386	42.70	ug/L	97
65) benzene	9.941	78	300496	54.56	ug/L	98
66) heptane	10.109	57	41851	48.86	ug/L	99
67) isopropyl acetate	9.858	43	136012	56.77	ug/L	100
68) 1,2-dichloroethane	9.952	62	95716	58.19	ug/L	99
69) trichloroethene	10.660	95	71275	53.53	ug/L	99
72) 2-nitropropane	11.425	41	17391	35.72	ug/L	94
74) methyl methacrylate	10.922	41	107374	57.02	ug/L	93
75) 1,2-dichloropropane	10.922	63	80495	55.91	ug/L	95
76) methylcyclohexane	10.875	83	123748	49.68	ug/L	98
77) dibromomethane	11.084	93	51861	61.23	ug/L	94
78) bromodichloromethane	11.210	83	102174	58.32	ug/L	99
79) cis-1,3-dichloropropene	11.672	75	116644	56.72	ug/L	94
81) 4-methyl-2-pentanone	11.766	58	32889	62.76	ug/L	99
82) toluene	12.033	92	169239	55.92	ug/L	99
83) 3-methyl-1-butanol	11.787	70	49696	1254.91	ug/L	96
84) trans-1,3-dichloropropene	12.233	75	107193	58.65	ug/L	99
85) ethyl methacrylate	12.227	69	86539	53.58	ug/L	99
86) 1,1,2-trichloroethane	12.448	83	57351	56.41	ug/L	98
87) 2-hexanone	12.621	58	24265	56.64	ug/L	97
89) tetrachloroethene	12.626	166	80218	48.33	ug/L	98
90) 1,3-dichloropropane	12.626	76	104947	55.48	ug/L	98
91) butyl acetate	12.694	56	46001	58.76	ug/L	96
92) 3,3-Dimethyl-1-Butanol	12.788	57	117562	628.33	ug/L	98
93) dibromochloromethane	12.893	129	86848	58.88	ug/L	99
94) 1,2-dibromoethane	13.045	107	72111	57.97	ug/L	98
95) n-Butyl Ether	13.423	57	287404	48.74	ug/L	99
96) chlorobenzene	13.507	112	193694	55.40	ug/L	100
97) 1,1,2,2-tetrachloroethane	13.564	131	93275	60.16	ug/L	99
98) ethylbenzene	13.559	91	313531	52.36	ug/L	99
99) m,p-xylene	13.664	106	245683	109.71	ug/L	99
100) o-xylene	14.083	106	128641	56.24	ug/L	100
101) styrene	14.094	104	203827	52.96	ug/L	99
103) bromoform	14.356	173	67500	58.65	ug/L	99
105) isopropylbenzene	14.424	105	332846	56.48	ug/L	99
107) bromobenzene	14.823	156	98900	56.86	ug/L	98
108) cyclohexanone	14.587	55	29910	190.61	ug/L	94
109) 1,1,2,2-tetrachloroethane	14.728	83	101377	59.45	ug/L	98
110) trans-1,4-dichloro-2-b...	14.770	53	18255	48.37	ug/L	97
111) 1,2,3-trichloropropane	14.802	110	23641	60.71	ug/L	90
112) n-propylbenzene	14.833	91	358137	55.82	ug/L	99
114) 2-chlorotoluene	14.980	126	87197	58.18	ug/L	98
115) 4-chlorotoluene	15.080	126	83732	57.54	ug/L	91
116) 1,3,5-trimethylbenzene	14.985	105	294268	58.68	ug/L	97
117) tert-butylbenzene	15.331	119	233643	56.54	ug/L	98
118) pentachloroethane	15.415	167	79829	70.84	ug/L	96
119) 1,2,4-trimethylbenzene	15.378	105	288522	57.31	ug/L	98
120) sec-butylbenzene	15.546	105	379751	57.51	ug/L	99
121) 1,3-dichlorobenzene	15.730	146	187717	58.05	ug/L	97
122) p-isopropyltoluene	15.662	119	330968	57.40	ug/L	99
123) 1,4-dichlorobenzene	15.814	146	192115	55.37	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
Data File : 3D112527.D
Acq On : 11 Aug 2015 4:30 pm
Operator : ximenac
Sample : jc1106-13msd
Misc : MS89468,V3D4822,5,,,1
ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 12 09:27:45 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.191	146	203019	58.23	ug/L	97
126) n-butylbenzene	16.065	92	165123	54.73	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.920	157	29079	57.07	ug/L	94
129) 1,3,5-trichlorobenzene	17.093	180	220653	59.61	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	232503	61.11	ug/L	97
131) hexachlorobutadiene	17.759	225	102558	54.45	ug/L	97
132) naphthalene	17.895	128	506611	62.86	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	246110	62.30	ug/L	98
134) hexachloroethane	16.443	119	70558	59.25	ug/L	98
135) Benzyl chloride	15.924	91	197569	58.32	ug/L	99

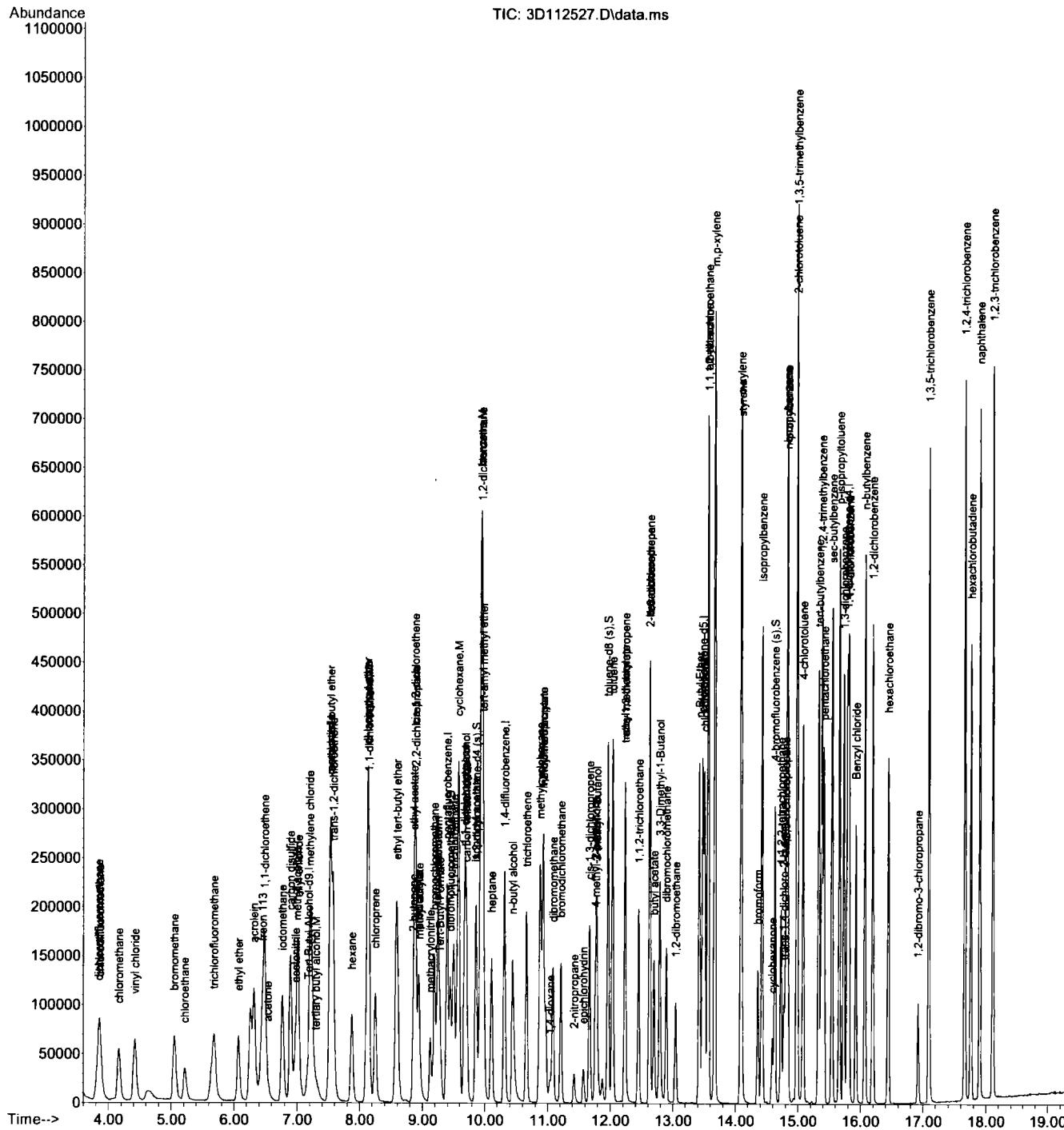
(#) = qualifier out of range (m) = manual integration (+) = signals summed

7.4.2
7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
Data File : 3D112527.D
Acq On : 11 Aug 2015 4:30 pm
Operator : ximenac
Sample : jc1106-13msd
Misc : MS89468,V3D4822,5,,,1
ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 12 09:27:45 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112566.D
 Acq On : 12 Aug 2015 10:44 am
 Operator : ximenac
 Sample : jc1072-1ms
 Misc : MS89457,V3D4823,S,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 12 14:30:35 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.178	65	99162	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	201673	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	227956	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	197596	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	122225	50.00	ug/L	0.00

System Monitoring Compounds

50) dibromofluoromethane (s)	9.438	113	85006	55.01	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 110.02%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	75958	53.07	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 106.14%	
80) toluene-d8 (s)	11.960	98	261661	52.41	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 104.82%	
106) 4-bromofluorobenzene (s)	14.629	95	94453	50.23	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.46%	

Target Compounds

					Qvalue
2) 1,4-dioxane	11.037	88	24075	1502.55	ug/L 92
3) tertiary butyl alcohol	7.299	59	53979	252.05	ug/L 97
8) chlorodifluoromethane	3.849	51	147116	53.74	ug/L 98
9) dichlorodifluoromethane	3.844	85	178550	48.71	ug/L 99
10) chloromethane	4.164	50	145660	49.98	ug/L 99
11) vinyl chloride	4.415	62	160381	54.87	ug/L 99
12) bromomethane	5.050	94	101640	53.82	ug/L 98
13) chloroethane	5.212	64	65743	53.10	ug/L 98
16) trichlorofluoromethane	5.674	101	176554	57.88	ug/L 98
18) ethyl ether	6.061	74	39802	52.82	ug/L 98
21) acrolein	6.318	56	143914	475.79	ug/L 99
22) 1,1-dichloroethene	6.481	61	121404	50.61	ug/L 99
23) acetone	6.533	43	23955	53.01	ug/L 83
24) allyl chloride	7.011	76	41232	55.34	ug/L 99
25) acetonitrile	6.969	40	85662	570.10	ug/L 82
27) iodomethane	6.764	142	190840	55.95	ug/L 97
28) iso-butyl alcohol	9.684	41	25910	516.05	ug/L 96
29) carbon disulfide	6.885	76	322550	53.17	ug/L 98
30) methylene chloride	7.199	84	98386	54.74	ug/L 100
31) methyl acetate	6.995	43	56126	55.91	ug/L 97
32) methyl tert butyl ether	7.530	73	284199	53.37	ug/L 99
33) trans-1,2-dichloroethene	7.572	61	116040	53.02	ug/L 98
34) di-isopropyl ether	8.122	45	280676	49.86	ug/L 92
35) ethyl tert-butyl ether	8.583	59	286134	53.95	ug/L 97
36) 2-butanone	8.846	72	7583	52.73	ug/L 98
37) 1,1-dichloroethane	8.138	63	143182	55.06	ug/L 97
38) chloroprene	8.248	53	94261	49.53	ug/L 96
39) acrylonitrile	7.524	53	152060	292.18	ug/L 96
40) vinyl acetate	8.122	86	11254	58.25	ug/L 86
41) ethyl acetate	8.856	45	8771	44.34	ug/L 93
42) 2,2-dichloropropane	8.877	77	150468	55.29	ug/L 99
43) cis-1,2-dichloroethene	8.872	96	87337	49.69	ug/L 99
44) propionitrile	8.935	54	109145	562.78	ug/L 95
45) methyl acrylate	8.940	55	72478	55.44	ug/L 97
46) bromochloromethane	9.186	128	47884	56.61	ug/L 96
47) tetrahydrofuran	9.228	42	24253	47.61	ug/L 99
48) chloroform	9.244	83	135615	54.09	ug/L 99
49) Tert-Butyl Formate	9.275	59	70116	48.04	ug/L # 94
52) freon 113	6.455	151	82726	56.01	ug/L 100
53) methacrylonitrile	9.123	41	36947	48.76	ug/L 98
54) 1,1,1-trichloroethane	9.496	97	142956	55.22	ug/L 97

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112566.D
 Acq On : 12 Aug 2015 10:44 am
 Operator : ximenac
 Sample : jc1072-1ms
 Misc : MS89457, V3D4823, 5, , , 1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 12 14:30:35 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
57) tert-amyl methyl ether	9.968	73	275041	51.89	ug/L	96
59) epichlorohydrin	11.572	57	30613	244.62	ug/L	98
60) n-butyl alcohol	10.439	56	98978	2659.32	ug/L	95
61) cyclohexane	9.569	84	128823	48.26	ug/L	97
62) carbon tetrachloride	9.700	117	129690	58.83	ug/L	98
63) 1,1-dichloropropene	9.674	75	86244	52.53	ug/L	99
64) hexane	7.870	57	73041	49.81	ug/L	99
65) benzene	9.936	78	281410	52.54	ug/L	97
66) heptane	10.104	57	42034	50.46	ug/L	94
67) isopropyl acetate	9.857	43	123295	52.92	ug/L	99
68) 1,2-dichloroethane	9.947	62	89897	56.19	ug/L	98
69) trichloroethene	10.654	95	68144	52.62	ug/L	100
72) 2-nitropropane	11.425	41	17382	36.71	ug/L	96
74) methyl methacrylate	10.922	41	97917	53.47	ug/L	92
75) 1,2-dichloropropane	10.922	63	74665	53.33	ug/L	98
76) methylcyclohexane	10.875	83	135116	55.77	ug/L	96
77) dibromomethane	11.079	93	49320	59.88	ug/L	94
78) bromodichloromethane	11.210	83	94853	55.67	ug/L	97
79) cis-1,3-dichloropropene	11.672	75	110594	55.30	ug/L	97
81) 4-methyl-2-pentanone	11.761	58	27476	53.91	ug/L	98
82) toluene	12.033	92	162686	55.27	ug/L	100
83) 3-methyl-1-butanol	11.787	70	40267	1045.53	ug/L	98
84) trans-1,3-dichloropropene	12.233	75	103407	58.18	ug/L	96
85) ethyl methacrylate	12.227	69	76000	48.38	ug/L	99
86) 1,1,2-trichloroethane	12.442	83	55076	55.70	ug/L	96
87) 2-hexanone	12.615	58	20285	48.69	ug/L	92
89) tetrachloroethene	12.621	166	77591	46.12	ug/L	96
90) 1,3-dichloropropane	12.626	76	99167	51.72	ug/L	97
91) butyl acetate	12.694	56	40733	51.34	ug/L	96
92) 3,3-Dimethyl-1-Butanol	12.783	57	80494	424.44	ug/L	99
93) dibromochloromethane	12.893	129	84074	56.23	ug/L	99
94) 1,2-dibromoethane	13.045	107	68014	53.95	ug/L	97
95) n-Butyl Ether	13.423	57	267174	44.70	ug/L	97
96) chlorobenzene	13.507	112	187493	52.91	ug/L	98
97) 1,1,1,2-tetrachloroethane	13.564	131	89179	56.75	ug/L	98
98) ethylbenzene	13.559	91	299257	49.31	ug/L	99
99) m,p-xylene	13.664	106	237455	104.62	ug/L	99
100) o-xylene	14.083	106	124912	53.88	ug/L	95
101) styrene	14.094	104	193974	49.73	ug/L	99
103) bromoform	14.356	173	64342	55.16	ug/L	99
105) isopropylbenzene	14.424	105	311571	50.80	ug/L	99
107) bromobenzene	14.823	156	96122	53.10	ug/L	98
108) cyclohexanone	14.587	55	26260	169.50	ug/L	97
109) 1,1,2,2-tetrachloroethane	14.728	83	98494	55.49	ug/L	99
110) trans-1,4-dichloro-2-b...	14.770	53	18140	46.18	ug/L	93
111) 1,2,3-trichloropropane	14.802	110	23275	57.42	ug/L	88
112) n-propylbenzene	14.833	91	340850	51.04	ug/L	99
114) 2-chlorotoluene	14.980	126	83855	53.76	ug/L	98
115) 4-chlorotoluene	15.080	126	82183	54.26	ug/L	92
116) 1,3,5-trimethylbenzene	14.985	105	275526	52.79	ug/L	99
117) tert-butylbenzene	15.331	119	209081	48.61	ug/L	98
118) pentachloroethane	15.410	167	75067	64.00	ug/L	97
119) 1,2,4-trimethylbenzene	15.378	105	274658	52.41	ug/L	98
120) sec-butylbenzene	15.546	105	342354	49.81	ug/L	99
121) 1,3-dichlorobenzene	15.730	146	184320	54.76	ug/L	99
122) p-isopropyltoluene	15.662	119	296579	49.42	ug/L	100
123) 1,4-dichlorobenzene	15.814	146	191548	53.04	ug/L	98
124) 1,2-dichlorobenzene	16.191	146	199943	55.10	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112566.D
 Acq On : 12 Aug 2015 10:44 am
 Operator : ximenac
 Sample : jc1072-1ms
 Misc : MS89457,V3D4823,5,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 12 14:30:35 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
126) n-butylbenzene	16.065	92	150910	48.06	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.920	157	27833	52.48	ug/L	95
129) 1,3,5-trichlorobenzene	17.093	180	202966	52.68	ug/L	99
130) 1,2,4-trichlorobenzene	17.659	180	211892	53.50	ug/L	97
131) hexachlorobutadiene	17.759	225	86841	44.30	ug/L	98
132) naphthalene	17.900	128	470470	56.08	ug/L	99
133) 1,2,3-trichlorobenzene	18.110	180	223450	54.34	ug/L	100
134) hexachloroethane	16.443	119	62277	50.24	ug/L	98
135) Benzyl chloride	15.924	91	201287	57.09	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

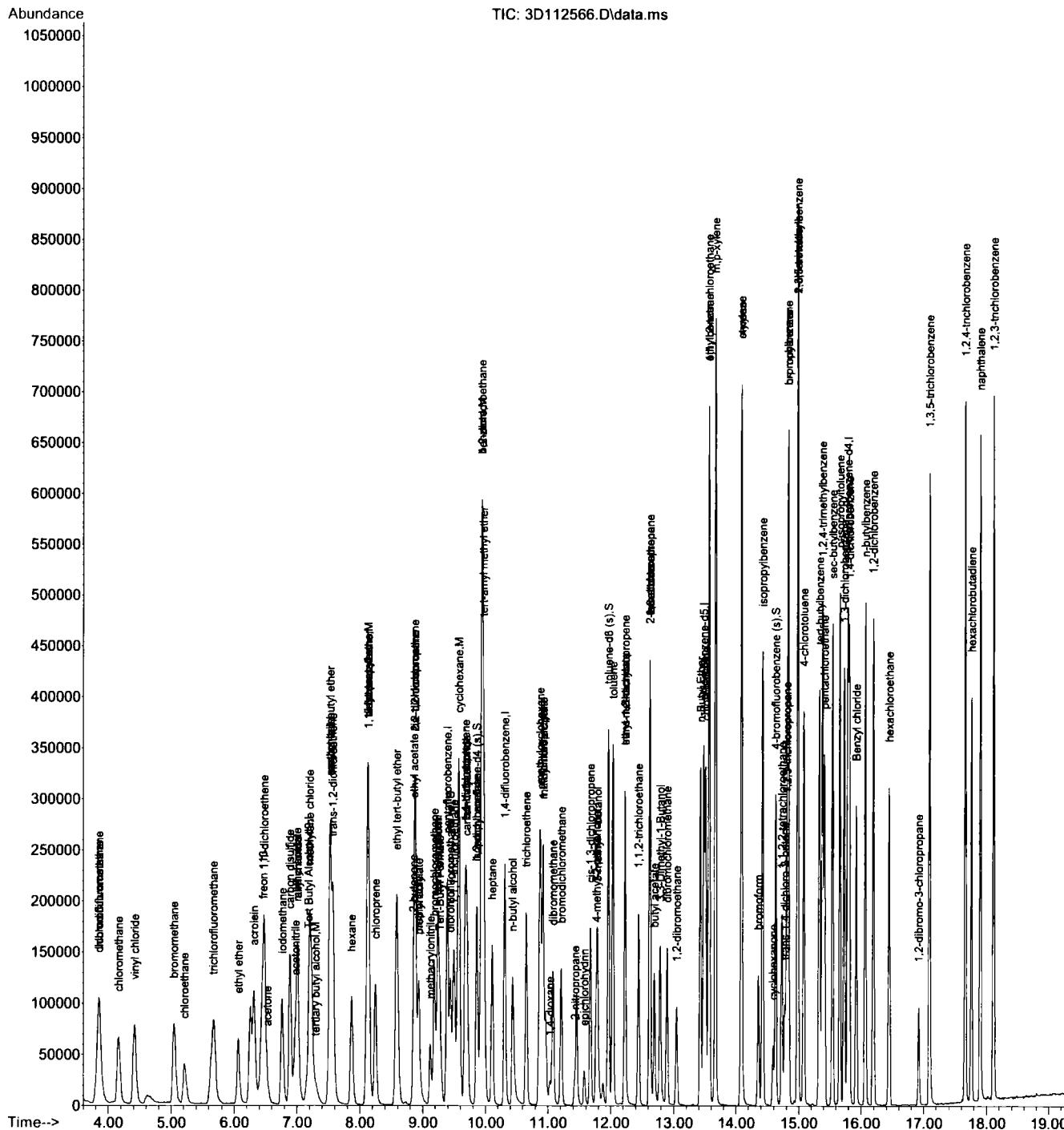
743

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Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
Data File : 3D112566.D
Acq On : 12 Aug 2015 10:44 am
Operator : ximenac
Sample : jc1072-1ms
Misc : MS89457,V3D4823,5,,,1
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 12 14:30:35 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112567.D
 Acq On : 12 Aug 2015 11:11 am
 Operator : ximenac
 Sample : jc1072-1msd
 Misc : MS89457,V3D4823,5,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 12 14:30:24 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.173	65	99080	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	197615	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	223819	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	186124	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	115672	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
50) dibromofluoromethane (s)	9.438	113	80888	53.42	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	= 106.84%		
51) 1,2-dichloroethane-d4 (s)	9.863	65	73491	52.40	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	= 104.80%		
80) toluene-d8 (s)	11.960	98	252659	51.54	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	= 103.08%		
106) 4-bromofluorobenzene (s)	14.629	95	88333	49.63	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	= 99.26%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-dioxane	11.032	88	25958	1621.41	ug/L	97
3) tertiary butyl alcohol	7.294	59	57556	268.98	ug/L	98
8) chlorodifluoromethane	3.844	51	147310	54.91	ug/L	98
9) dichlorodifluoromethane	3.844	85	179706	50.03	ug/L	98
10) chloromethane	4.158	50	142921	50.04	ug/L	98
11) vinyl chloride	4.410	62	160854	56.16	ug/L	100
12) bromomethane	5.044	94	99067	53.53	ug/L	98
13) chloroethane	5.207	64	62789	51.75	ug/L	98
16) trichlorofluoromethane	5.668	101	172450	57.69	ug/L	100
18) ethyl ether	6.061	74	40756	55.20	ug/L	97
21) acrolein	6.318	56	146250	493.44	ug/L	96
22) 1,1-dichloroethene	6.481	61	118926	50.60	ug/L	97
23) acetone	6.533	43	23270	52.55	ug/L	99
24) allyl chloride	7.010	76	39021	53.45	ug/L	98
25) acetonitrile	6.963	40	85366	579.79	ug/L	81
27) iodomethane	6.764	142	187707	56.16	ug/L	97
28) iso-butyl alcohol	9.684	41	26954	547.87	ug/L	98
29) carbon disulfide	6.885	76	313369	52.72	ug/L	99
30) methylene chloride	7.199	84	96406	54.74	ug/L	95
31) methyl acetate	6.995	43	57358	58.31	ug/L	97
32) methyl tert butyl ether	7.529	73	288749	55.34	ug/L	99
33) trans-1,2-dichloroethene	7.571	61	114081	53.20	ug/L	97
34) di-isopropyl ether	8.117	45	285559	51.77	ug/L	98
35) ethyl tert-butyl ether	8.589	59	288200	55.46	ug/L	98
36) 2-butanone	8.840	72	8338	59.17	ug/L	87
37) 1,1-dichloroethane	8.138	63	142043	55.75	ug/L	99
38) chloroprene	8.243	53	94122	50.47	ug/L	98
39) acrylonitrile	7.524	53	155258	304.45	ug/L	98
40) vinyl acetate	8.122	86	11687	61.74	ug/L	93
41) ethyl acetate	8.856	45	9189	47.40	ug/L	79
42) 2,2-dichloropropane	8.877	77	147270	55.23	ug/L	99
43) cis-1,2-dichloroethene	8.877	96	86902	50.46	ug/L	97
44) propionitrile	8.935	54	112888	594.04	ug/L	95
45) methyl acrylate	8.940	55	75590	59.01	ug/L	99
46) bromochloromethane	9.186	128	47774	57.64	ug/L	94
47) tetrahydrofuran	9.223	42	24254	48.59	ug/L	96
48) chloroform	9.239	83	133924	54.51	ug/L	99
49) Tert-Butyl Formate	9.275	59	60765	42.48	ug/L #	96
52) freon 113	6.449	151	81369	56.23	ug/L	96
53) methacrylonitrile	9.123	41	38950	52.46	ug/L	98
54) 1,1,1-trichloroethane	9.501	97	138107	54.45	ug/L	97

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112567.D
 Acq On : 12 Aug 2015 11:11 am
 Operator : ximenac
 Sample : jc1072-1msd
 Misc : MS89457,V3D4823,5,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 12 14:30:24 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
57) tert-amyl methyl ether	9.968	73	279001	53.71	ug/L	96
59) epichlorohydrin	11.577	57	30931	251.73	ug/L	99
60) n-butyl alcohol	10.439	56	105489	2886.64	ug/L	100
61) cyclohexane	9.569	84	123316	47.05	ug/L	99
62) carbon tetrachloride	9.700	117	125204	57.84	ug/L	99
63) 1,1-dichloropropene	9.674	75	85899	53.28	ug/L	99
64) hexane	7.870	57	72730	50.51	ug/L	97
65) benzene	9.936	78	279672	53.18	ug/L	99
66) heptane	10.104	57	43314	52.95	ug/L	98
67) isopropyl acetate	9.857	43	127326	55.66	ug/L	99
68) 1,2-dichloroethane	9.952	62	91054	57.97	ug/L	96
69) trichloroethene	10.654	95	68351	53.76	ug/L	98
72) 2-nitropropane	11.425	41	15980	34.38	ug/L	96
74) methyl methacrylate	10.922	41	98843	54.97	ug/L	92
75) 1,2-dichloropropane	10.922	63	75451	54.88	ug/L	100
76) methylcyclohexane	10.875	83	133784	56.24	ug/L	96
77) dibromomethane	11.079	93	48991	60.58	ug/L	94
78) bromodichloromethane	11.210	83	95182	56.90	ug/L	98
79) cis-1,3-dichloropropene	11.672	75	110784	56.41	ug/L	99
81) 4-methyl-2-pentanone	11.761	58	28776	57.50	ug/L	97
82) toluene	12.033	92	161940	56.03	ug/L	96
83) 3-methyl-1-butanol	11.782	70	41813	1105.74	ug/L	99
84) trans-1,3-dichloropropene	12.233	75	104000	59.59	ug/L	99
85) ethyl methacrylate	12.227	69	79952	51.84	ug/L	98
86) 1,1,2-trichloroethane	12.442	83	56502	58.20	ug/L	94
87) 2-hexanone	12.621	58	21228	51.89	ug/L	96
89) tetrachloroethene	12.626	166	76923	48.54	ug/L	99
90) 1,3-dichloropropane	12.626	76	102200	56.59	ug/L	98
91) butyl acetate	12.694	56	41854	56.00	ug/L	95
92) 3,3-Dimethyl-1-Butanol	12.788	57	87612	490.45	ug/L	98
93) dibromochloromethane	12.893	129	84180	59.78	ug/L	98
94) 1,2-dibromoethane	13.045	107	68658	57.81	ug/L	95
95) n-Butyl Ether	13.423	57	265039	47.07	ug/L	98
96) chlorobenzene	13.507	112	186305	55.82	ug/L	98
97) 1,1,1,2-tetrachloroethane	13.564	131	87992	59.44	ug/L	97
98) ethylbenzene	13.559	91	293412	51.32	ug/L	100
99) m,p-xylene	13.664	106	231748	108.40	ug/L	99
100) o-xylene	14.083	106	119960	54.93	ug/L	98
101) styrene	14.094	104	191532	52.13	ug/L	99
103) bromoform	14.361	173	65014	59.17	ug/L	98
105) isopropylbenzene	14.424	105	300078	51.70	ug/L	99
107) bromobenzene	14.823	156	93104	54.34	ug/L	98
108) cyclohexanone	14.587	55	27452	181.41	ug/L	94
109) 1,1,2,2-tetrachloroethane	14.728	83	99067	58.97	ug/L	97
110) trans-1,4-dichloro-2-b...	14.770	53	18555	49.91	ug/L	91
111) 1,2,3-trichloropropane	14.802	110	23055	60.10	ug/L	92
112) n-propylbenzene	14.833	91	325766	51.55	ug/L	99
114) 2-chlorotoluene	14.980	126	81181	54.99	ug/L	94
115) 4-chlorotoluene	15.079	126	77405	54.00	ug/L	97
116) 1,3,5-trimethylbenzene	14.985	105	260889	52.82	ug/L	98
117) tert-butylbenzene	15.331	119	205120	50.39	ug/L	99
118) pentachloroethane	15.410	167	73021	65.78	ug/L	97
119) 1,2,4-trimethylbenzene	15.378	105	262120	52.85	ug/L	100
120) sec-butylbenzene	15.546	105	330061	50.75	ug/L	99
121) 1,3-dichlorobenzene	15.730	146	176192	55.31	ug/L	98
122) p-isopropyltoluene	15.661	119	282594	49.76	ug/L	99
123) 1,4-dichlorobenzene	15.813	146	182598	53.43	ug/L	99
124) 1,2-dichlorobenzene	16.191	146	191926	55.89	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
Data File : 3D112567.D
Acq On : 12 Aug 2015 11:11 am
Operator : ximenac
Sample : jc1072-1msd
Misc : MS89457,V3D4823.5,,,1
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 12 14:30:24 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration

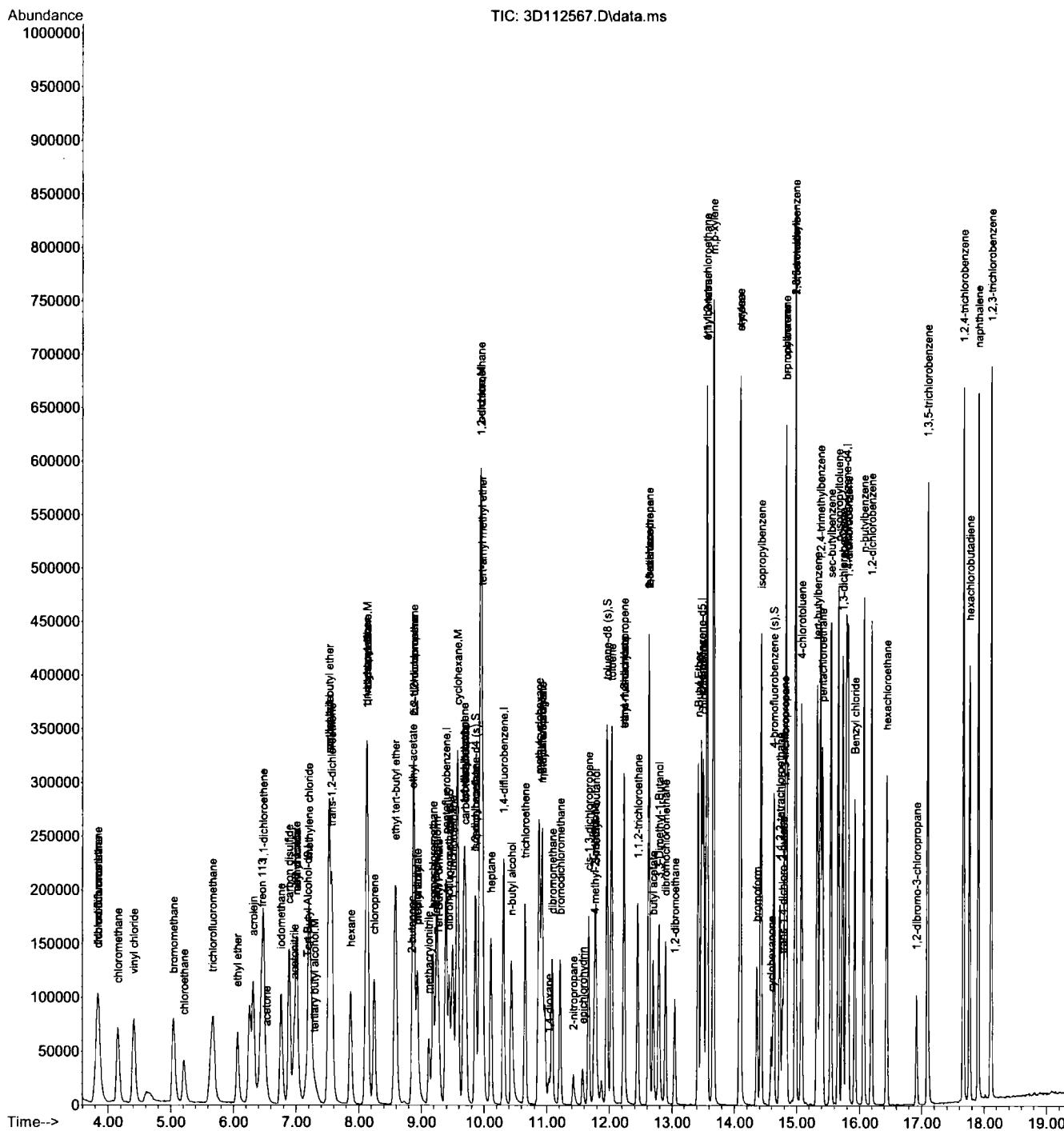
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
126) n-butylbenzene	16.065	92	144293	48.55	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.920	157	28816	57.42	ug/L	97
129) 1,3,5-trichlorobenzene	17.093	180	194015	53.21	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	208872	55.73	ug/L	97
131) hexachlorobutadiene	17.759	225	89643	48.32	ug/L	98
132) naphthalene	17.895	128	472061	59.46	ug/L	99
133) 1,2,3-trichlorobenzene	18.110	180	220243	56.60	ug/L	99
134) hexachloroethane	16.437	119	60758	51.79	ug/L	98
135) Benzyl chloride	15.924	91	199350	59.74	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
Data File : 3D112567.D
Acq On : 12 Aug 2015 11:11 am
Operator : ximenac
Sample : jc1072-1msd
Misc : MS89457,V3D4823,5,,,1
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 12 14:30:24 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
 Data File : 3D112580.D
 Acq On : 12 Aug 2015 5:04 pm
 Operator : ximenac
 Sample : jc1106-9ms
 Misc : MS89468,V3D4824,5,,,1
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 13 10:55:18 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.183	65	118942	500.00	ug/L	0.00
4) pentafluorobenzene	9.396	168	243673	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	269432	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	220893	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	129859	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	99592	53.34	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	=	106.68%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	84836	49.05	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	=	98.10%	
80) toluene-d8 (s)	11.960	98	300148	50.86	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	=	101.72%	
106) 4-bromofluorobenzene (s)	14.629	95	102824	51.46	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	=	102.92%	
Target Compounds						
					Qvalue	
2) 1,4-dioxane	11.037	88	29277	1523.35	ug/L	98
3) tertiary butyl alcohol	7.299	59	68870	268.10	ug/L	98
8) chlorodifluoromethane	3.859	51	149713	45.26	ug/L	100
9) dichlorodifluoromethane	3.849	85	176433	39.84	ug/L	99
10) chloromethane	4.169	50	156778	44.52	ug/L	99
11) vinyl chloride	4.420	62	173369	49.09	ug/L	100
12) bromomethane	5.050	94	111056	48.67	ug/L	98
13) chloroethane	5.217	64	69034	46.15	ug/L	98
16) trichlorofluoromethane	5.684	101	178449	48.42	ug/L	96
18) ethyl ether	6.072	74	48221	52.96	ug/L	96
21) acrolein	6.324	56	170834	467.44	ug/L	98
22) 1,1-dichloroethene	6.486	61	144097	49.72	ug/L	94
23) acetone	6.544	43	28154	51.57	ug/L	91
24) allyl chloride	7.010	76	46587	51.75	ug/L	89
25) acetonitrile	6.984	40	99990	550.75	ug/L	91
27) iodomethane	6.769	142	234986	57.02	ug/L	97
28) iso-butyl alcohol	9.690	41	33327	549.37	ug/L	92
29) carbon disulfide	6.895	76	367546	50.15	ug/L	99
30) methylene chloride	7.204	84	117471	54.09	ug/L	97
31) methyl acetate	7.000	43	68074	56.12	ug/L	97
32) methyl tert butyl ether	7.529	73	344903	53.61	ug/L	99
33) trans-1,2-dichloroethene	7.577	61	134850	50.99	ug/L	97
34) di-isopropyl ether	8.122	45	336711	49.51	ug/L	92
35) ethyl tert-butyl ether	8.589	59	346892	54.13	ug/L	98
36) 2-butanone	8.856	72	9539	54.90	ug/L	91
37) 1,1-dichloroethane	8.143	63	168581	53.66	ug/L	99
38) chloroprene	8.248	53	110288	47.96	ug/L	96
39) acrylonitrile	7.529	53	185809	295.49	ug/L	94
40) vinyl acetate	8.132	86	13945	59.74	ug/L	75
41) ethyl acetate	8.866	45	10421	43.60	ug/L #	1
42) 2,2-dichloropropane	8.877	77	167652	50.99	ug/L	99
43) cis-1,2-dichloroethene	8.877	96	114685	54.01	ug/L	96
44) propionitrile	8.940	54	136437	582.25	ug/L	93
45) methyl acrylate	8.940	55	87706	55.52	ug/L	98
46) bromochloromethane	9.186	128	58993	57.73	ug/L	96
47) tetrahydrofuran	9.228	42	28663	46.57	ug/L	95
48) chloroform	9.244	83	161039	53.16	ug/L	98
49) Tert-Butyl Formate	9.270	59	77644	44.02	ug/L #	97
52) freon 113	6.455	151	92379	51.77	ug/L	100
53) methacrylonitrile	9.123	41	44740	48.87	ug/L	96

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
 Data File : 3D112580.D
 Acq On : 12 Aug 2015 5:04 pm
 Operator : ximenac
 Sample : jc1106-9ms
 Misc : MS89468,V3D4824,5,,,1
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 13 10:55:18 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
54) 1,1,1-trichloroethane	9.496	97	245937	78.63	ug/L	98
57) tert-amyl methyl ether	9.973	73	333272	52.04	ug/L	99
59) epichlorohydrin	11.577	57	37667	254.66	ug/L	93
60) n-butyl alcohol	10.439	56	134425	3055.72	ug/L	95
61) cyclohexane	9.569	84	146766	46.52	ug/L	97
62) carbon tetrachloride	9.705	117	141640	54.36	ug/L	98
63) 1,1-dichloropropene	9.674	75	99863	51.46	ug/L	97
64) hexane	7.876	57	85248	49.18	ug/L	99
65) benzene	9.936	78	329016	51.97	ug/L	99
66) heptane	10.109	57	48794	49.56	ug/L	95
67) isopropyl acetate	9.857	43	159239	57.82	ug/L	99
68) 1,2-dichloroethane	9.952	62	102119	54.01	ug/L	99
69) trichloroethene	10.660	95	80069	52.31	ug/L	96
72) 2-nitropropane	11.425	41	17612	31.47	ug/L	99
74) methyl methacrylate	10.922	41	116858	53.99	ug/L	92
75) 1,2-dichloropropane	10.922	63	86601	52.33	ug/L	100
76) methylcyclohexane	10.875	83	156528	54.67	ug/L	95
77) dibromomethane	11.079	93	56175	57.70	ug/L	94
78) bromodichloromethane	11.210	83	108951	54.11	ug/L	99
79) cis-1,3-dichloropropene	11.671	75	126627	53.57	ug/L	95
81) 4-methyl-2-pentanone	11.761	58	36431	60.47	ug/L	98
82) toluene	12.033	92	186346	53.56	ug/L	99
83) 3-methyl-1-butanol	11.782	70	57708	1267.73	ug/L	94
84) trans-1,3-dichloropropene	12.232	75	115504	54.98	ug/L	96
85) ethyl methacrylate	12.227	69	99339	53.50	ug/L	94
86) 1,1,2-trichloroethane	12.442	83	63873	54.65	ug/L	93
87) 2-hexanone	12.620	58	27687	56.22	ug/L	96
89) tetrachloroethene	12.626	166	88433	47.02	ug/L	99
90) 1,3-dichloropropane	12.631	76	114159	53.26	ug/L	96
91) butyl acetate	12.694	56	52922	59.66	ug/L	93
92) 3,3-Dimethyl-1-Butanol	12.788	57	137227	647.28	ug/L	98
93) dibromochloromethane	12.893	129	95996	57.44	ug/L	99
94) 1,2-dibromoethane	13.045	107	79380	56.32	ug/L	95
95) n-Butyl Ether	13.423	57	316783	47.41	ug/L	99
96) chlorobenzene	13.507	112	214270	54.09	ug/L	98
97) 1,1,1,2-tetrachloroethane	13.564	131	104109	59.26	ug/L	99
98) ethylbenzene	13.559	91	345227	50.88	ug/L	98
99) m,p-xylene	13.664	106	271633	107.05	ug/L	98
100) o-xylene	14.083	106	146051	56.35	ug/L	97
101) styrene	14.094	104	223149	51.17	ug/L	99
103) bromoform	14.361	173	74117	56.84	ug/L	98
105) isopropylbenzene	14.424	105	369534	56.71	ug/L	99
107) bromobenzene	14.823	156	108038	56.17	ug/L	95
108) cyclohexanone	14.587	55	32221	187.12	ug/L	96
109) 1,1,2,2-tetrachloroethane	14.728	83	113463	60.16	ug/L	99
110) trans-1,4-dichloro-2-b...	14.770	53	13312	31.89	ug/L	91
111) 1,2,3-trichloropropane	14.802	110	25724	59.73	ug/L	93
112) n-propylbenzene	14.833	91	389070	54.84	ug/L	100
114) 2-chlorotoluene	14.980	126	93938	56.68	ug/L	96
115) 4-chlorotoluene	15.079	126	91952	57.14	ug/L	89
116) 1,3,5-trimethylbenzene	14.985	105	321149	57.91	ug/L	99
117) tert-butylbenzene	15.331	119	252891	55.34	ug/L	99
118) pentachloroethane	15.410	167	86846	69.69	ug/L	98
119) 1,2,4-trimethylbenzene	15.378	105	320802	57.62	ug/L	98
120) sec-butylbenzene	15.546	105	413153	56.58	ug/L	99
121) 1,3-dichlorobenzene	15.730	146	203280	56.84	ug/L	99
122) p-isopropyltoluene	15.661	119	354869	55.66	ug/L	98
123) 1,4-dichlorobenzene	15.813	146	207439	54.07	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
 Data File : 3D112580.D
 Acq On : 12 Aug 2015 5:04 pm
 Operator : ximenac
 Sample : jc1106-9ms
 Misc : MS89468,V3D4824,5,,,1
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 13 10:55:18 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.191	146	218588	56.70	ug/L	97
126) n-butylbenzene	16.065	92	174682	52.36	ug/L	98
128) 1,2-dibromo-3-chloropr...	16.920	157	31729	56.31	ug/L	97
129) 1,3,5-trichlorobenzene	17.093	180	232551	56.81	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	244660	58.15	ug/L	97
131) hexachlorobutadiene	17.759	225	99495	47.77	ug/L	98
132) naphthalene	17.895	128	558007	62.61	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	257406	58.92	ug/L	97
134) hexachloroethane	16.443	119	75009	56.95	ug/L	98
135) Benzyl chloride	15.924	91	225235	60.13	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\

Data File : 3D112580.D

Acq On : 12 Aug 2015 5:04 pm

Operator : ximenac

Sample : jc1106-9ms

Misc : MS89468, V3D4824.5, . . . 1

ALS Vial : 19 Sample Multiplier: 1

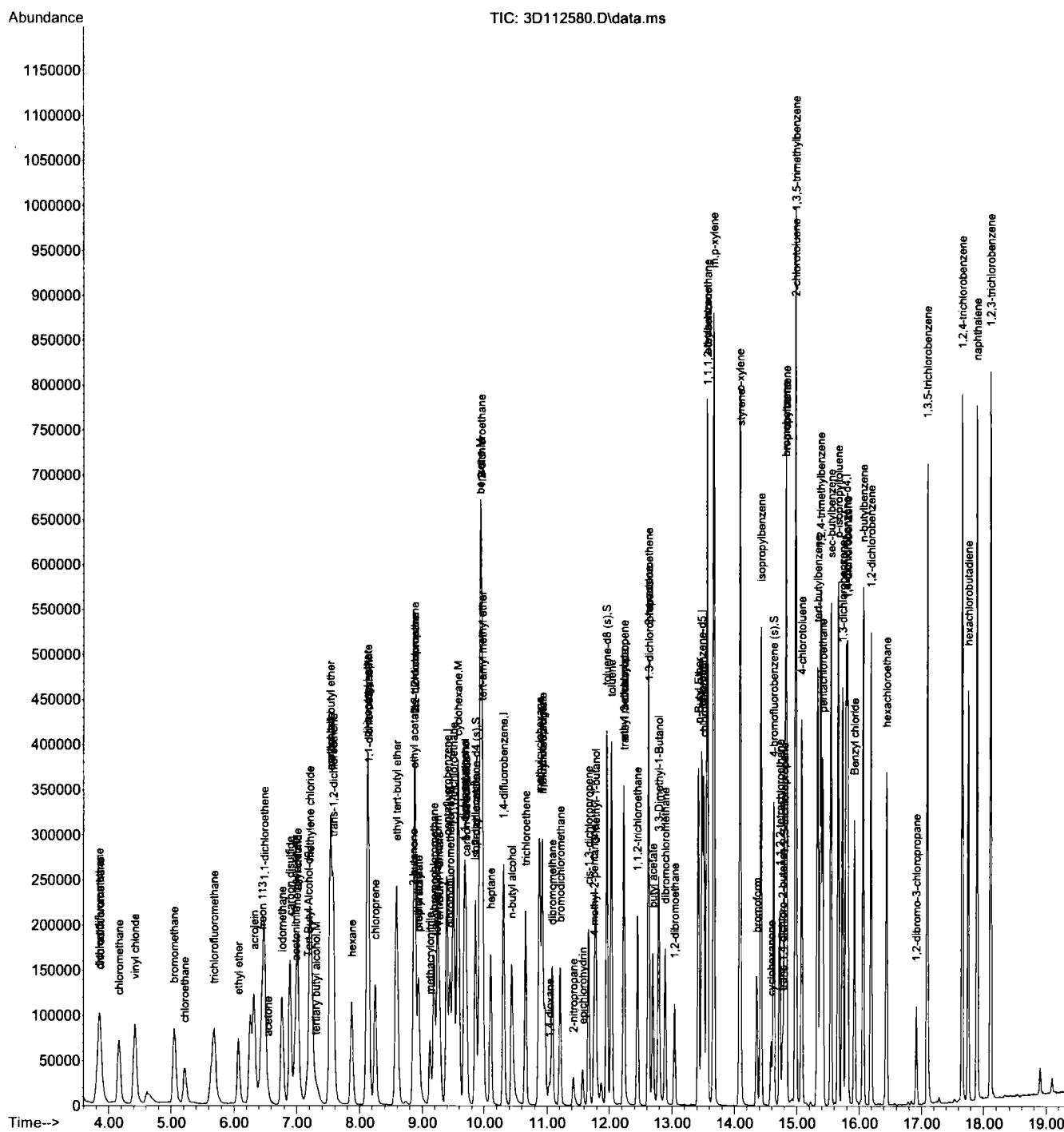
Quant Time: Aug 13 10:55:18 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Method : Method SW846 8260B. ZB 624 60m x 0.25mm x 1.4 um

Last Update : Thu Aug 06 11:10:54 2015

Qbase update : Thu Aug 08 11:10:54
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
 Data File : 3D112581.D
 Acq On : 12 Aug 2015 5:31 pm
 Operator : ximenac
 Sample : jc1106-9msd
 Misc : MS89468,V3D4824,5,,,1
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 13 10:55:33 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.178	65	120440	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	245404	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.313	114	271506	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	223373	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	132733	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	98227	52.24	ug/L	0.00
Spiked Amount 50.000 Range 76 - 120			Recovery	=	104.48%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	84093	48.28	ug/L	0.00
Spiked Amount 50.000 Range 73 - 122			Recovery	=	96.56%	
80) toluene-d8 (s)	11.960	98	304159	51.15	ug/L	0.00
Spiked Amount 50.000 Range 84 - 119			Recovery	=	102.30%	
106) 4-bromofluorobenzene (s)	14.628	95	103529	50.69	ug/L	0.00
Spiked Amount 50.000 Range 78 - 117			Recovery	=	101.38%	
Target Compounds						
2) 1,4-dioxane	11.037	88	31259	1606.25	ug/L	98
3) tertiary butyl alcohol	7.283	59	71369	274.38	ug/L	93
8) chlorodifluoromethane	3.859	51	150846	45.28	ug/L	98
9) dichlorodifluoromethane	3.844	85	180947	40.57	ug/L	99
10) chloromethane	4.163	50	153746	43.35	ug/L	99
11) vinyl chloride	4.420	62	173771	48.85	ug/L	98
12) bromomethane	5.049	94	110789	48.21	ug/L	97
13) chloroethane	5.212	64	68041	45.16	ug/L	98
16) trichlorofluoromethane	5.679	101	180751	48.69	ug/L	99
18) ethyl ether	6.072	74	48214	52.58	ug/L	93
21) acrolein	6.318	56	165424	449.45	ug/L	99
22) 1,1-dichloroethene	6.486	61	142451	48.80	ug/L	95
23) acetone	6.538	43	26747	48.64	ug/L	93
24) allyl chloride	7.016	76	48676	53.69	ug/L	93
25) acetonitrile	6.979	40	98817	540.45	ug/L	96
27) iodomethane	6.764	142	235196	56.67	ug/L	98
28) iso-butyl alcohol	9.684	41	33465	547.75	ug/L	93
29) carbon disulfide	6.895	76	367504	49.79	ug/L	98
30) methylene chloride	7.204	84	116382	53.21	ug/L	95
31) methyl acetate	7.000	43	65984	54.02	ug/L	96
32) methyl tert butyl ether	7.529	73	341097	52.64	ug/L	98
33) trans-1,2-dichloroethene	7.571	61	133109	49.98	ug/L	99
34) di-isopropyl ether	8.122	45	335679	49.01	ug/L	93
35) ethyl tert-butyl ether	8.583	59	345296	53.51	ug/L	98
36) 2-butanone	8.845	72	9056	51.75	ug/L	100
37) 1,1-dichloroethane	8.138	63	166248	52.54	ug/L	99
38) chloroprene	8.248	53	107384	46.37	ug/L	96
39) acrylonitrile	7.529	53	178307	281.56	ug/L	95
40) vinyl acetate	8.127	86	13928	59.25	ug/L	88
41) ethyl acetate	8.856	45	11684	48.54	ug/L	61
42) 2,2-dichloropropane	8.882	77	164890	49.80	ug/L	98
43) cis-1,2-dichloroethene	8.877	96	112305	52.51	ug/L	96
44) propionitrile	8.935	54	131291	556.34	ug/L	92
45) methyl acrylate	8.940	55	85954	54.03	ug/L	99
46) bromochloromethane	9.186	128	58126	56.48	ug/L	95
47) tetrahydrofuran	9.223	42	27569	44.47	ug/L	93
48) chloroform	9.244	83	158266	51.87	ug/L	100
49) Tert-Butyl Formate	9.275	59	60646	34.14	ug/L	# 97
52) freon 113	6.455	151	92868	51.67	ug/L	96
53) methacrylonitrile	9.123	41	44948	48.75	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\

Data File : 3D112581.D

Acq On : 12 Aug 2015 5:31 pm

Operator : ximenac

Sample : jc1106-9msd

Misc : MS89468,V3D4824,5,,,1

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 13 10:55:33 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Aug 06 11:10:54 2015

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
54) 1,1,1-trichloroethane	9.501	97	247216	78.48	ug/L	97
57) tert-amyl methyl ether	9.967	73	332323	51.52	ug/L	98
59) epichlorohydrin	11.572	57	36351	243.88	ug/L	98
60) n-butyl alcohol	10.439	56	134782	3040.43	ug/L	96
61) cyclohexane	9.574	84	146344	46.03	ug/L	96
62) carbon tetrachloride	9.705	117	141526	53.90	ug/L	98
63) 1,1-dichloropropene	9.674	75	97659	49.94	ug/L	99
64) hexane	7.870	57	83558	47.84	ug/L	98
65) benzene	9.936	78	329060	51.58	ug/L	98
66) heptane	10.109	57	48456	48.84	ug/L	97
67) isopropyl acetate	9.857	43	154136	55.54	ug/L	98
68) 1,2-dichloroethane	9.952	62	101170	53.10	ug/L	99
69) trichloroethene	10.654	95	80528	52.21	ug/L	99
72) 2-nitropropane	11.425	41	17196	30.49	ug/L	94
74) methyl methacrylate	10.922	41	115705	53.05	ug/L	92
75) 1,2-dichloropropane	10.922	63	86939	52.13	ug/L	98
76) methylcyclohexane	10.874	83	154924	53.69	ug/L	96
77) dibromomethane	11.079	93	55998	57.08	ug/L	95
78) bromodichloromethane	11.210	83	108595	53.52	ug/L	97
79) cis-1,3-dichloropropene	11.671	75	127624	53.58	ug/L	98
81) 4-methyl-2-pentanone	11.761	58	36247	59.71	ug/L	99
82) toluene	12.033	92	189899	54.17	ug/L	99
83) 3-methyl-1-butanol	11.787	70	57545	1254.49	ug/L	94
84) trans-1,3-dichloropropene	12.232	75	118174	55.82	ug/L	93
85) ethyl methacrylate	12.227	69	98254	52.51	ug/L	97
86) 1,1,2-trichloroethane	12.447	83	64781	55.01	ug/L	97
87) 2-hexanone	12.620	58	27699	55.82	ug/L	96
89) tetrachloroethene	12.626	166	90263	47.46	ug/L	97
90) 1,3-dichloropropane	12.626	76	117108	54.03	ug/L	97
91) butyl acetate	12.694	56	52366	58.38	ug/L	92
92) 3,3-Dimethyl-1-Butanol	12.788	57	136313	635.83	ug/L	98
93) dibromochloromethane	12.893	129	95866	56.72	ug/L	99
94) 1,2-dibromoethane	13.045	107	79829	56.01	ug/L	96
95) n-Butyl Ether	13.423	57	314784	46.59	ug/L	98
96) chlorobenzene	13.506	112	219227	54.73	ug/L	97
97) 1,1,1,2-tetrachloroethane	13.564	131	103486	58.25	ug/L	99
98) ethylbenzene	13.559	91	346330	50.48	ug/L	99
99) m,p-xylene	13.664	106	270932	105.59	ug/L	97
100) o-xylene	14.083	106	144745	55.23	ug/L	92
101) styrene	14.094	104	223372	50.66	ug/L	99
103) bromoform	14.356	173	73545	55.77	ug/L	100
105) isopropylbenzene	14.424	105	369051	55.41	ug/L	99
107) bromobenzene	14.822	156	108505	55.19	ug/L	97
108) cyclohexanone	14.587	55	32872	186.88	ug/L	96
109) 1,1,2,2-tetrachloroethane	14.728	83	112065	58.14	ug/L	99
110) trans-1,4-dichloro-2-b...	14.770	53	14496	33.98	ug/L	99
111) 1,2,3-trichloropropane	14.802	110	25718	58.42	ug/L	90
112) n-propylbenzene	14.833	91	389072	53.65	ug/L	99
114) 2-chlorotoluene	14.980	126	94247	55.64	ug/L	95
115) 4-chlorotoluene	15.079	126	91518	55.64	ug/L	89
116) 1,3,5-trimethylbenzene	14.985	105	320652	56.57	ug/L	98
117) tert-butylbenzene	15.331	119	260725	55.82	ug/L	96
118) pentachloroethane	15.415	167	86425	67.85	ug/L	99
119) 1,2,4-trimethylbenzene	15.378	105	315291	55.40	ug/L	98
120) sec-butylbenzene	15.546	105	415688	55.70	ug/L	100
121) 1,3-dichlorobenzene	15.735	146	204439	55.93	ug/L	99
122) p-isopropyltoluene	15.661	119	353525	54.25	ug/L	99
123) 1,4-dichlorobenzene	15.813	146	208322	53.12	ug/L	100

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
Data File : 3D112581.D
Acq On : 12 Aug 2015 5:31 pm
Operator : ximenac
Sample : jc1106-9msd
Misc : MS89468,V3D4824,5,,,1
ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 13 10:55:33 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration

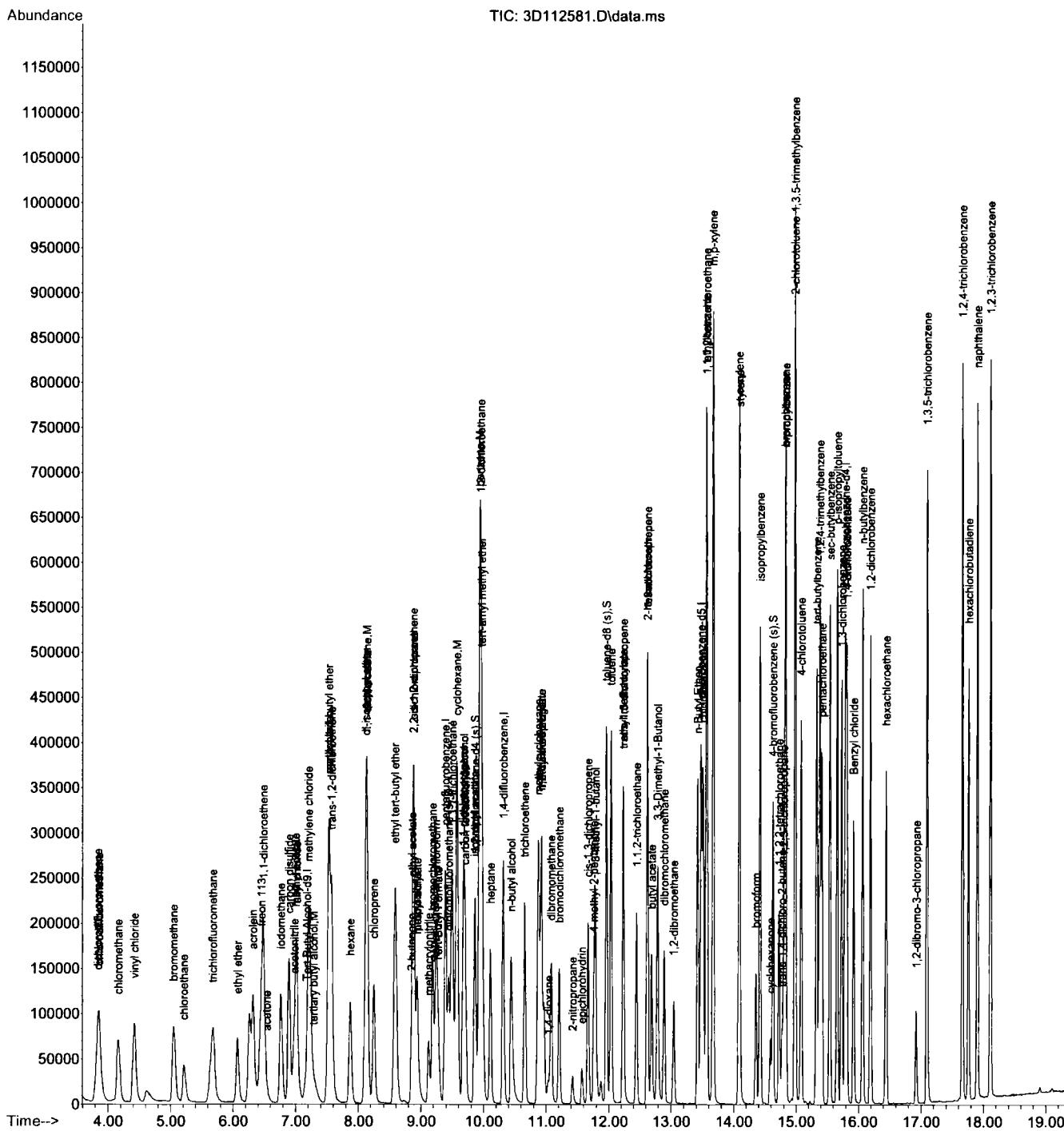
Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
124) 1,2-dichlorobenzene	16.191	146	221300	56.16	ug/L	98
126) n-butylbenzene	16.065	92	173046	50.75	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.925	157	31134	54.06	ug/L	98
129) 1,3,5-trichlorobenzene	17.093	180	235947	56.39	ug/L	99
130) 1,2,4-trichlorobenzene	17.659	180	251677	58.52	ug/L	96
131) hexachlorobutadiene	17.759	225	104875	49.26	ug/L	99
132) naphthalene	17.900	128	557097	61.15	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	267604	59.93	ug/L	99
134) hexachloroethane	16.443	119	74829	55.59	ug/L	98
135) Benzyl chloride	15.924	91	222934	58.22	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4823-4824\
Data File : 3D112581.D
Acq On : 12 Aug 2015 5:31 pm
Operator : ximenac
Sample : jc1106-9msd
Misc : MS89468,V3D4824,5,,,1
ALS Vial : 20 Sample Multiplier: 1

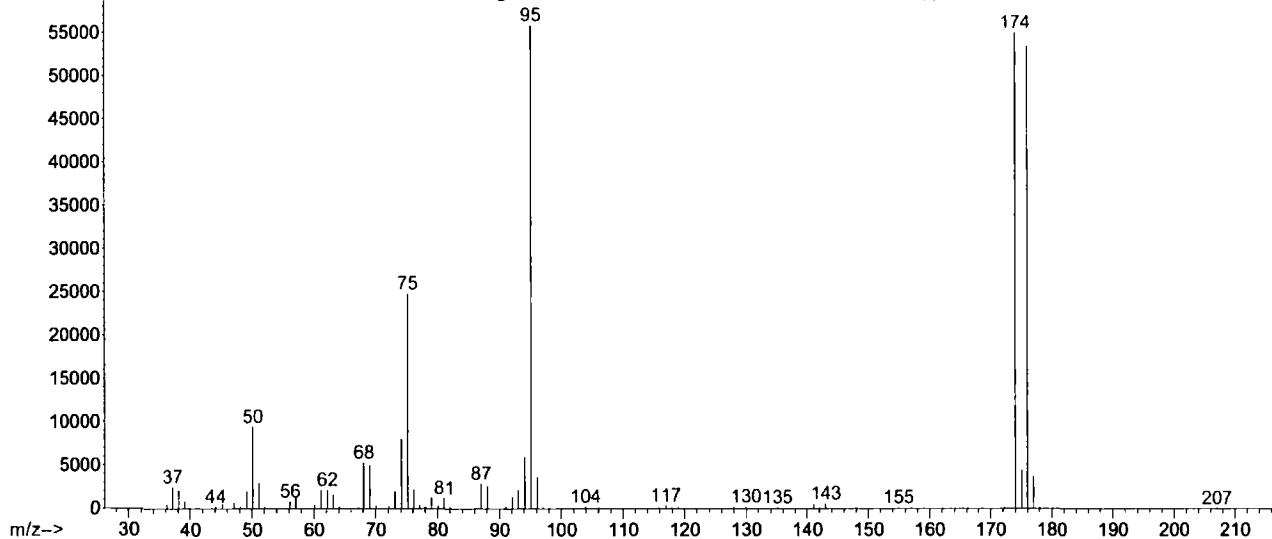
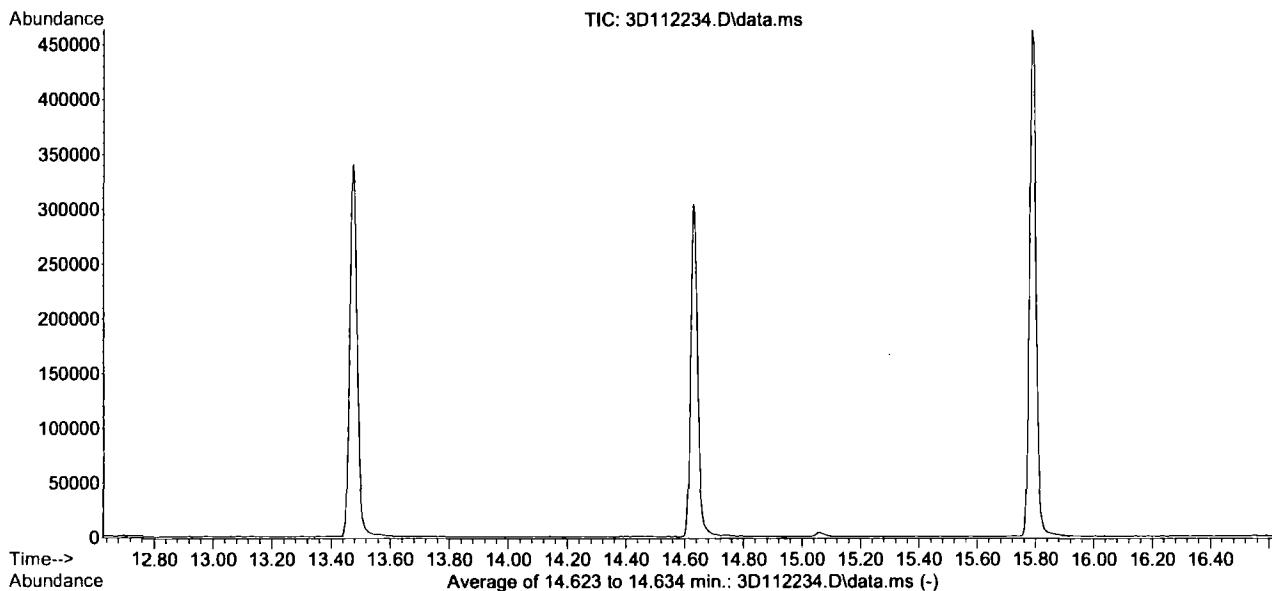
Quant Time: Aug 13 10:55:33 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 B260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



SW-846 Method 8260

Data File : C:\msdchem\1\DATA\3D112234.D Vial: 19
 Acq On : 29 Jul 2015 7:16 pm Operator: ximenac
 Sample : bfb Inst : MS3D
 Misc : MS88759,V3D4810,5,,,1 Multiplr: 1.00
 MS Integration Params: rteint.p

Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)
 Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um



AutoFind: Scans 2105, 2106, 2107; Background Corrected with Scan 2096

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	16.9	9449	PASS
75	95	30	60	44.4	24824	PASS
95	95	100	100	100.0	55917	PASS
96	95	5	9	6.5	3622	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	98.4	55002	PASS
175	174	5	9	8.0	4400	PASS
176	174	95	101	97.3	53496	PASS
177	176	5	9	6.9	3680	PASS

Average of 14.623 to 14.634 min.: 3D112234.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.05	493	50.05	9449	64.00	176	77.70	93
37.05	2471	51.05	2986	67.10	126	77.95	253
38.05	2152	52.00	148	68.00	5281	78.90	1247
39.00	879	55.00	90	69.00	4980	80.00	329
39.85	75	56.05	800	70.00	366	80.95	1205
41.00	3	57.05	1312	72.00	256	81.90	232
44.00	189	58.00	27	73.00	2007	85.95	63
45.05	534	60.05	448	74.00	8028	86.95	2893
47.05	714	61.05	2114	75.05	24824	88.00	2603
47.95	270	62.05	2139	76.05	2203	90.95	262
49.05	1993	63.00	1570	77.00	380	92.00	1308

Average of 14.623 to 14.634 min.: 3D112234.D\data.ms

bfb

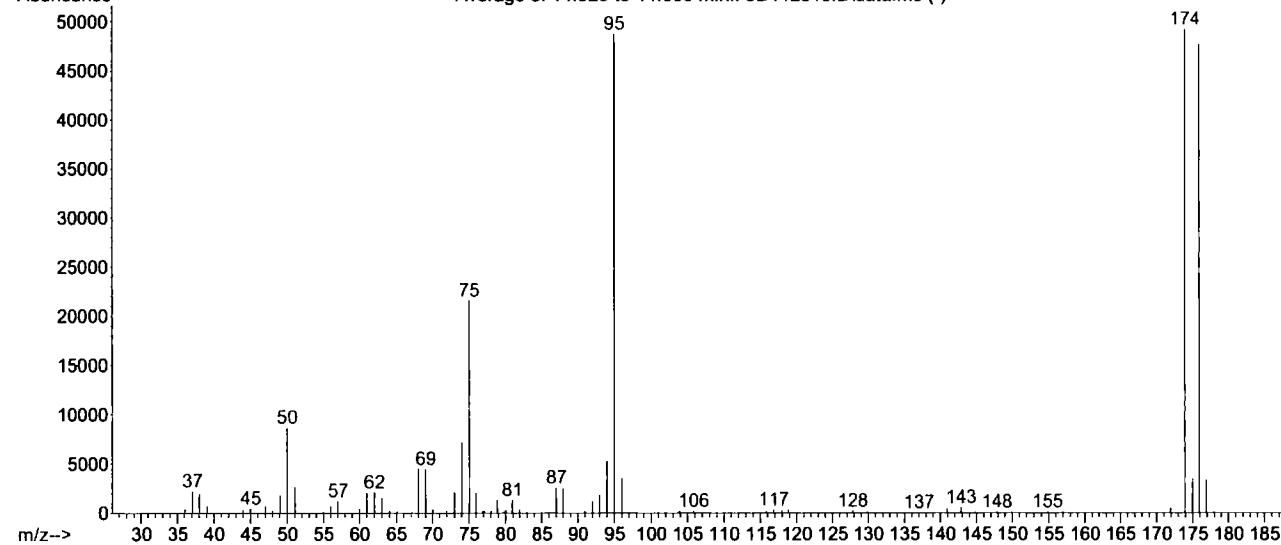
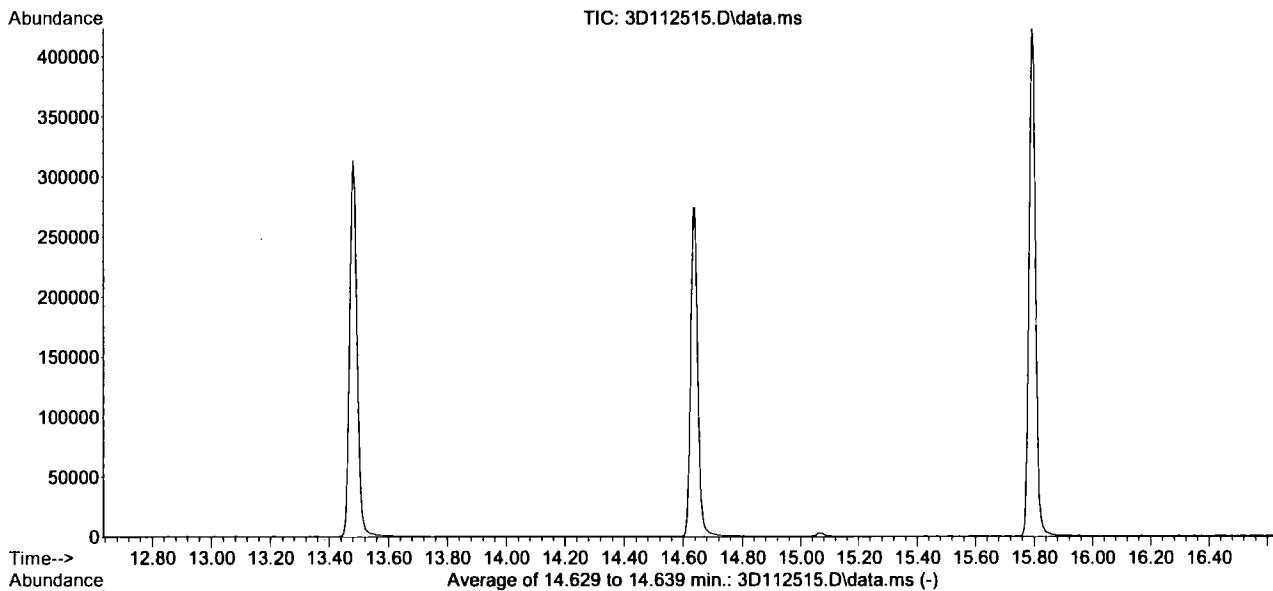
Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
93.00	2121	116.90	341	145.90	39	177.85	116
94.05	5943	117.85	145	147.90	149	206.95	14
95.00	55917	118.90	246	154.85	123		
96.00	3622	127.90	167	156.95	68		
97.00	131	128.90	35	159.00	36		
103.85	210	129.85	199	171.60	26		
104.80	27	134.95	117	172.15	123		
105.00	67	136.90	43	173.90	55002		
105.95	184	140.95	440	175.00	4400		
114.90	29	141.85	74	175.90	53496		
115.85	160	142.90	533	176.90	3680		

SW-846 Method 8260

Data File : C:\msdchem\1\DATA\3D112515.D Vial: 3
 Acq On : 11 Aug 2015 9:53 am Operator: ximenac
 Sample : bfb Inst : MS3D
 Misc : MS89341,V3D4822,5,,,,1 Multiplr: 1.00
 MS Integration Params: rteint.p

Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)
 Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 μ m



AutoFind: Scans 2106, 2107, 2108; Background Corrected with Scan 2098

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	17.8	8665	PASS
75	95	30	60	44.2	21568	PASS
95	95	100	100	100.0	48757	PASS
96	95	5	9	7.1	3479	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	100.9	49194	PASS
175	174	5	9	7.2	3521	PASS
176	174	95	101	96.9	47664	PASS
177	176	5	9	6.9	3271	PASS

Average of 14.629 to 14.639 min.: 3D112515.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.05	361	51.05	2687	67.10	135	78.05	194
37.05	2209	52.05	72	68.00	4455	78.90	1278
38.05	1937	54.95	113	69.00	4470	79.80	144
39.10	741	56.00	655	70.05	363	80.05	270
39.95	16	57.00	1190	71.95	248	80.95	1302
44.00	316	60.00	405	73.00	2109	81.95	303
45.00	454	61.00	2063	74.05	7123	86.00	26
47.00	696	62.00	2092	75.05	21568	87.00	2573
48.00	248	63.00	1519	76.00	2035	88.00	2413
49.05	1771	64.05	224	76.90	228	90.95	152
50.00	8665	65.05	163	77.10	165	92.00	1168

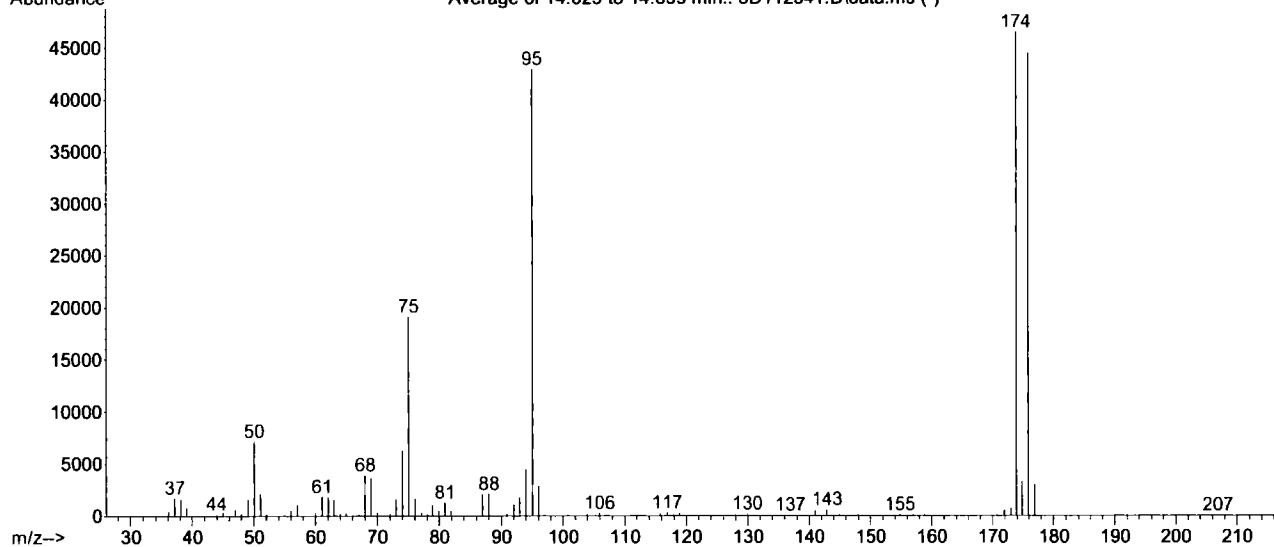
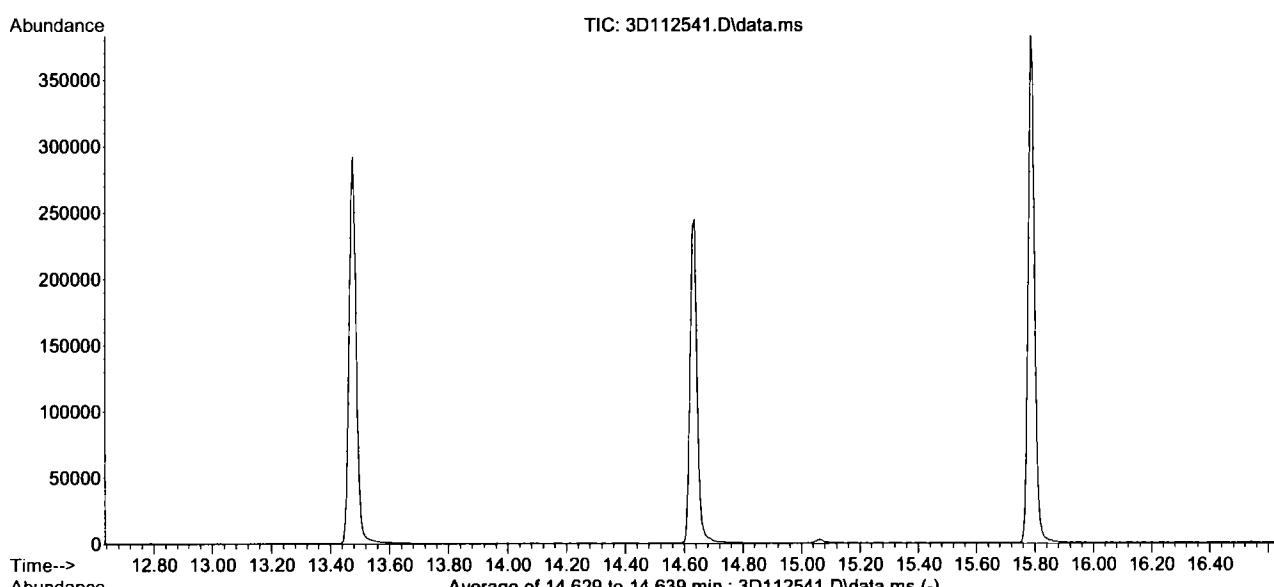
Average of 14.629 to 14.639 min.: 3D112515.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
93.00	1778	116.85	317	142.85	570	175.90	47664
94.00	5278	117.95	180	144.85	99	176.90	3271
95.00	48757	118.90	281	147.85	106	177.95	68
96.05	3479	127.85	197	149.90	36		
96.90	54	129.00	36	154.85	127		
103.80	79	129.90	155	156.80	31		
103.95	126	130.80	30	157.00	67		
104.90	28	135.00	30	170.90	26		
105.85	195	137.00	30	171.95	513		
114.90	25	140.95	431	173.90	49194		
115.80	188	142.10	32	174.95	3521		

SW-846 Method 8260
 Data File : C:\msdchem\1\DATA\3D\V3D4822\3D112541.D Vial: 34
 Acq On : 11 Aug 2015 10:50 pm Operator: ximenac
 Sample : bfb Inst : MS3D
 Misc : MS89468,V3D4823,5,,,1 Multiplr: 1.00
 MS Integration Params: rteint.p
 Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)
 Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um



AutoFind: Scans 2106, 2107, 2108; Background Corrected with Scan 2097

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	16.6	7142	PASS
75	95	30	60	44.5	19101	PASS
95	95	100	100	100.0	42928	PASS
96	95	5	9	6.6	2834	PASS
173	174	0.00	2	1.4	632	PASS
174	95	50	120	108.2	46456	PASS
175	174	5	9	7.1	3302	PASS
176	174	95	101	95.8	44520	PASS
177	176	5	9	6.6	2956	PASS

Average of 14.629 to 14.639 min.: 3D112541.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.05	362	51.10	2113	67.05	147	78.90	1091
37.10	1686	52.05	110	68.00	3901	79.95	446
38.10	1599	54.95	79	69.00	3608	80.90	1258
39.05	745	56.00	484	70.00	302	81.90	387
39.95	1	57.00	1032	72.00	203	87.00	2076
43.95	137	60.00	351	73.00	1577	88.00	2111
45.00	321	61.00	1831	74.00	6264	90.95	176
47.00	571	62.05	1765	75.00	19101	92.00	1103
47.95	244	63.00	1475	76.05	1637	93.00	1810
49.05	1557	63.95	182	77.05	287	94.00	4462
50.05	7142	64.95	267	78.05	194	95.00	42928

Average of 14.629 to 14.639 min.: 3D112541.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
96.05	2834	134.95	92	170.70	54		
96.95	92	136.90	62	171.90	548		
103.95	193	140.95	444	173.00	632		
105.95	221	142.90	589	173.90	46456		
115.80	187	144.90	80	174.90	3302		
116.90	289	146.00	28	175.90	44520		
117.90	159	147.90	121	176.95	2956		
118.95	232	154.85	88	206.90	49		
127.95	149	155.90	37				
128.90	63	156.95	60				
129.90	182	158.85	60				

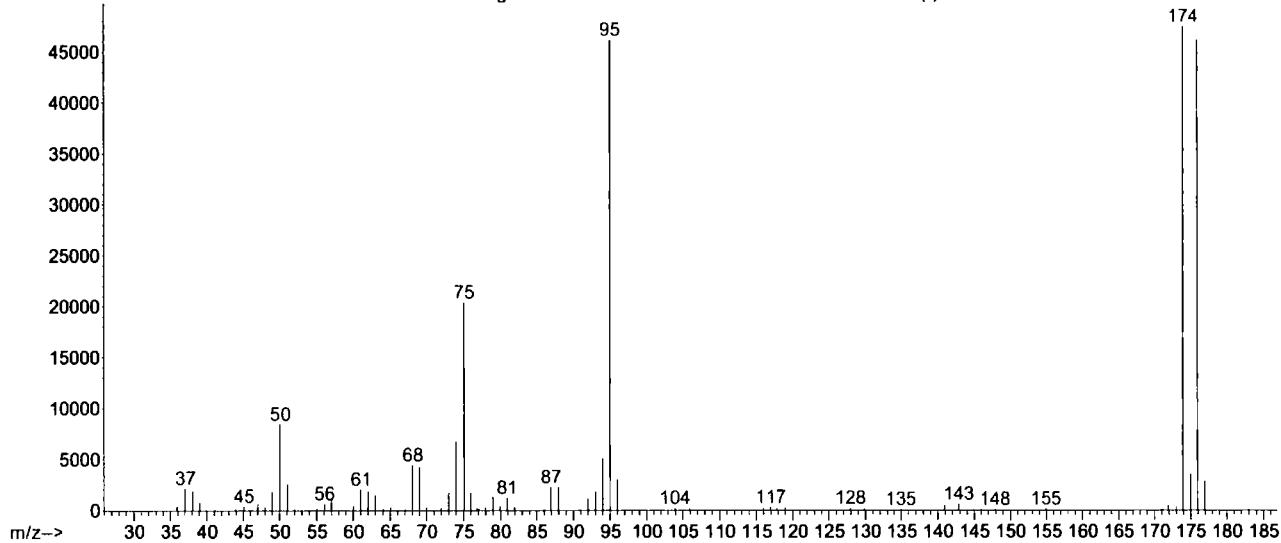
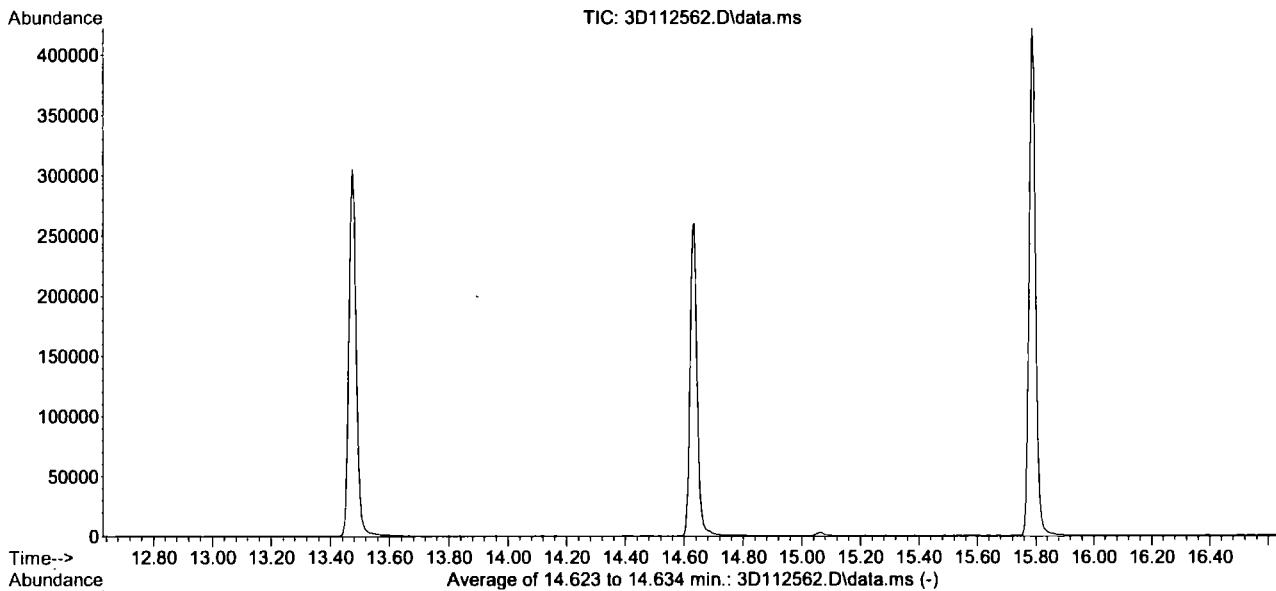
7.5.3

7

SW-846 Method 8260

Data File : C:\msdchem\1\DATA\3D112562.D Vial: 1
 Acq On : 12 Aug 2015 8:16 am Operator: ximenac
 Sample : bfb Inst : MS3D
 Misc : MS89457,V3D4824,5,,,1 Multiplr: 1.00
 MS Integration Params: rteint.p

Method : C:\msdchem\1\METHODS\M3D4810.M (RTE Integrator)
 Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um



AutoFind: Scans 2105, 2106, 2107; Background Corrected with Scan 2097

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	18.2	8422	PASS
75	95	30	60	44.1	20373	PASS
95	95	100	100	100.0	46160	PASS
96	95	5	9	6.4	2967	PASS
173	174	0.00	2	0.8	385	PASS
174	95	50	120	102.8	47440	PASS
175	174	5	9	7.5	3561	PASS
176	174	95	101	97.4	46184	PASS
177	176	5	9	6.2	2874	PASS

Average of 14.623 to 14.634 min.: 3D112562.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
35.95	390	49.00	1789	64.05	179	76.85	148
37.05	2149	50.05	8422	65.10	283	77.05	152
38.10	1907	51.05	2554	67.05	100	77.95	303
39.05	779	52.05	117	68.05	4438	78.95	1375
39.95	56	55.05	115	69.00	4241	79.90	388
41.10	30	56.05	614	69.95	305	80.90	1243
43.95	113	57.00	1014	71.95	210	81.95	344
45.05	364	60.05	434	73.00	1754	85.90	34
46.00	27	61.00	2028	74.00	6704	86.10	35
47.00	677	62.00	1884	75.05	20373	86.95	2252
48.05	259	63.00	1467	76.00	1689	87.95	2249

Average of 14.623 to 14.634 min.: 3D112562.D\data.ms

bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
90.80	47	114.80	35	142.85	588	174.95	3561
90.95	103	115.95	142	144.00	29	175.90	46184
92.00	1155	116.95	291	144.90	159	176.95	2874
93.00	1866	117.85	145	145.95	68		
94.00	5090	118.95	227	147.90	107		
95.00	46160	127.95	206	154.90	112		
96.00	2967	129.85	170	170.95	55		
96.80	34	130.90	30	171.20	47		
97.00	34	134.90	86	171.90	472		
103.90	202	136.80	31	173.00	385		
105.90	194	140.90	463	173.90	47440		

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4810-4810\
 Data File : 3d112235.d
 Acq On : 29 Jul 2015 8:15 pm
 Operator : ximenac
 Sample : IC4810-0.2
 Misc : MS88759,V3D4810.5,,,1
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 06 13:48:03 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.178	65	110638	500.00	ug/L	0.00
4) pentafluorobenzene	9.401	168	213329	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.319	114	234555	50.00	ug/L	0.00
88) chlorobenzene-d5	13.480	117	191394	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.793	152	121226	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)
50) dibromofluoromethane (s)	9.448	113	80221	43.48	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery =	86.96%		
51) 1,2-dichloroethane-d4 (s)	9.873	65	75864	39.83	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery =	79.66%		
80) toluene-d8 (s)	11.970	98	258418	51.74	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery =	103.48%		
106) 4-bromofluorobenzene (s)	14.634	95	92882	45.28	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	90.56%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) chloromethane	4.158	50	847	0.27	ug/L	# 48
11) vinyl chloride	4.415	62	488	0.14	ug/L	# 50
12) bromomethane	5.060	94	417	0.24	ug/L	# 60
27) iodomethane	6.775	142	469	0.11	ug/L	88
29) carbon disulfide	6.895	76	1035	0.11	ug/L	71
30) methylene chloride	7.225	84	326	0.13	ug/L	# 33
32) methyl tert butyl ether	7.556	73	1049	0.15	ug/L	# 46
33) trans-1,2-dichloroethene	7.598	61	361	0.11	ug/L	# 83
34) di-isopropyl ether	8.143	45	1282	0.16	ug/L	# 46
35) ethyl tert-butyl ether	8.589	59	829	0.11	ug/L	75
37) 1,1-dichloroethane	8.164	63	407	0.11	ug/L	# 49
38) chloroprene	8.264	53	303	0.11	ug/L	37
39) acrylonitrile	7.566	53	381	0.54	ug/L	# 60
42) 2,2-dichloropropane	8.903	77	631	0.18	ug/L	56
43) cis-1,2-dichloroethene	8.893	96	302	0.14	ug/L	# 41
48) chloroform	9.260	83	463	0.13	ug/L	59
57) tert-amyl methyl ether	9.973	73	1114	0.15	ug/L	52
62) carbon tetrachloride	9.726	117	374	0.15	ug/L	# 66
64) hexane	7.881	57	287	0.14	ug/L	87
65) benzene	9.947	78	1004	0.15	ug/L	92
69) trichloroethene	10.675	95	245	0.17	ug/L	75
76) methylcyclohexane	10.901	83	328	0.11	ug/L	# 21
78) bromodichloromethane	11.226	83	294	0.15	ug/L	# 27
79) cis-1,3-dichloropropene	11.692	75	369	0.16	ug/L	# 61
82) toluene	12.060	92	453	0.13	ug/L	52
84) trans-1,3-dichloropropene	12.259	75	275	0.14	ug/L	# 67
96) chlorobenzene	13.512	112	521	0.13	ug/L	65
98) ethylbenzene	13.575	91	1026	0.15	ug/L	83
99) m,p-xylene	13.690	106	607	0.24	ug/L	72
100) o-xylene	14.104	106	262	0.11	ug/L	# 42
105) isopropylbenzene	14.429	105	890	0.13	ug/L	84
109) 1,1,2,2-tetrachloroethane	14.739	83	307	0.15	ug/L	92
112) n-propylbenzene	14.854	91	963	0.12	ug/L	96
115) 4-chlorotoluene	15.111	126	227	0.14	ug/L	# 30
116) 1,3,5-trimethylbenzene	14.996	105	732	0.12	ug/L	90
117) tert-butylbenzene	15.336	119	608	0.12	ug/L	99
119) 1,2,4-trimethylbenzene	15.394	105	838	0.14	ug/L	95
120) sec-butylbenzene	15.557	105	788	0.10	ug/L	79
121) 1,3-dichlorobenzene	15.745	146	489	0.13	ug/L	91
122) p-isopropyltoluene	15.682	119	816	0.12	ug/L	68
123) 1,4-dichlorobenzene	15.819	146	642	0.16	ug/L	91

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\

Data File : 3d112235.d

Acq On : 29 Jul 2015 8:15 pm

Operator : ximenac

Sample : IC4810-0.2

Misc : MS88759,V3D4810,5,,,1

ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 06 13:48:03 2015

Quant Method : C:\msdchem\1\METHODS\M3D4810.M

Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um

QLast Update : Thu Jul 30 08:37:34 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.207	146	557	0.14	ug/L	99
129) 1,3,5-trichlorobenzene	17.114	180	477	0.11	ug/L #	62
130) 1,2,4-trichlorobenzene	17.675	180	630	0.15	ug/L	96
132) naphthalene	17.916	128	1654	0.17	ug/L	94
133) 1,2,3-trichlorobenzene	18.120	180	628	0.14	ug/L	96
135) Benzyl chloride	15.945	91	651	0.18	ug/L #	61

(#) = qualifier out of range (m) = manual integration (+) = signals summed

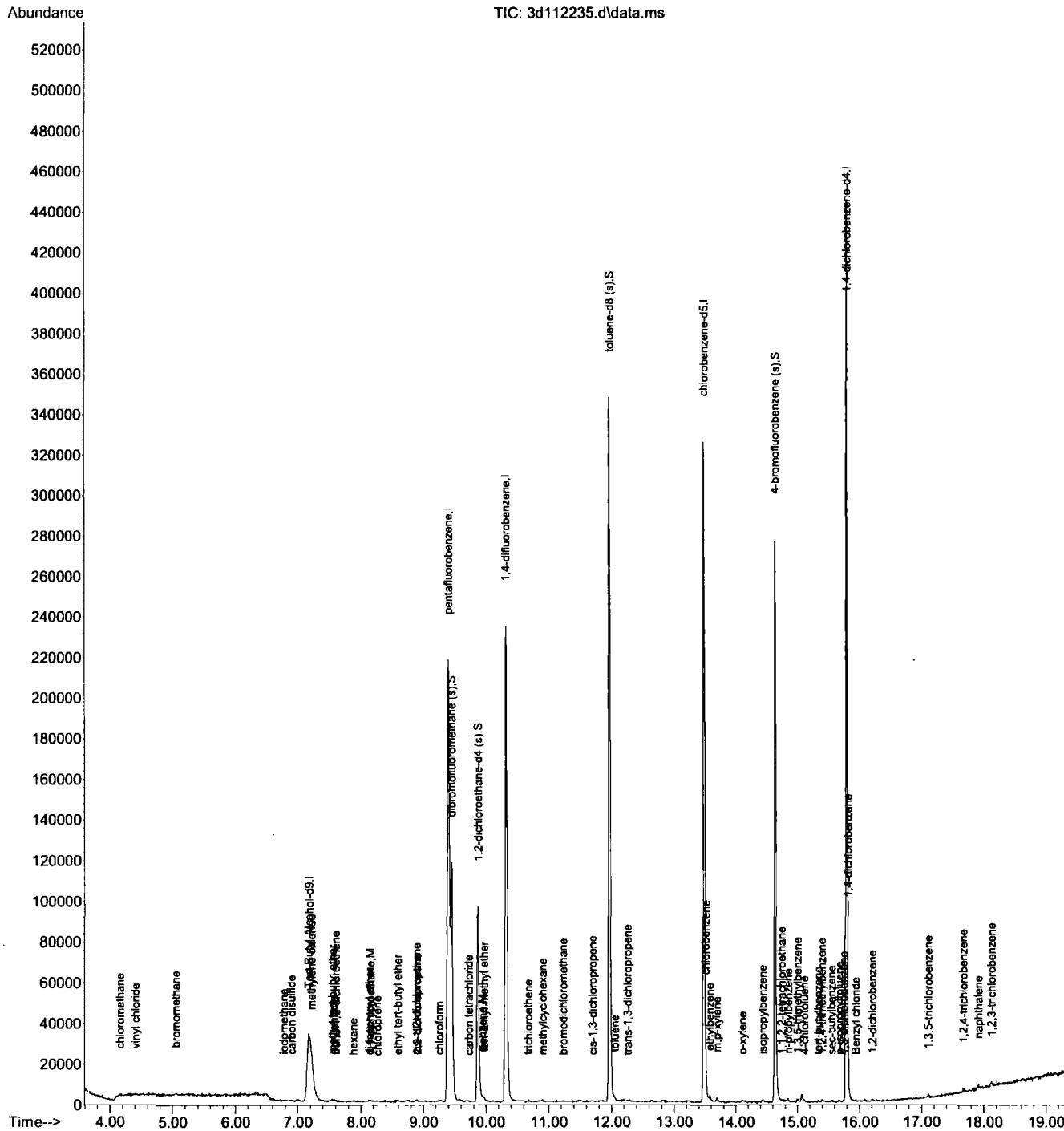
7.61

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Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112235.d
 Acq On : 29 Jul 2015 8:15 pm
 Operator : ximenac
 Sample : IC4810-0.2
 Misc : MS88759, V3D4810, 5, , , 1
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 06 13:48:03 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112236.d
 Acq On : 29 Jul 2015 8:42 pm
 Operator : ximenac
 Sample : IC4810-0.5
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Aug 06 13:48:29 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.189	65	96167	500.00	ug/L	0.01
4) pentafluorobenzene	9.401	168	202817	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.319	114	226728	50.00	ug/L	0.00
88) chlorobenzene-d5	13.480	117	188072	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.792	152	123641	50.00	ug/L	0.00

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System Monitoring Compounds

50) dibromofluoromethane (s)	9.454	113	78153	44.55	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	89.10%
51) 1,2-dichloroethane-d4 (s)	9.873	65	73280	40.46	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	80.92%
80) toluene-d8 (s)	11.970	98	246952	51.15	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	102.30%
106) 4-bromofluorobenzene (s)	14.634	95	94678	45.25	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.50%

Target Compounds

				Qvalue
10) chloromethane	4.153	50	1710	0.57 ug/L 84
11) vinyl chloride	4.410	62	1538	0.47 ug/L 69
12) bromomethane	5.055	94	1144	0.70 ug/L # 52
13) chloroethane	5.233	64	796	0.65 ug/L 70
16) trichlorofluoromethane	5.694	101	1436	0.43 ug/L 87
21) acrolein	6.339	56	1567	3.98 ug/L 68
22) 1,1-dichloroethene	6.486	61	849	0.24 ug/L 75
24) allyl chloride	7.010	76	257	0.25 ug/L # 13
27) iodomethane	6.780	142	1508	0.36 ug/L 97
29) carbon disulfide	6.911	76	2697	0.31 ug/L 73
30) methylene chloride	7.225	84	917	0.38 ug/L # 51
32) methyl tert butyl ether	7.535	73	2637	0.40 ug/L 96
33) trans-1,2-dichloroethene	7.592	61	1028	0.33 ug/L # 63
34) di-isopropyl ether	8.132	45	3114	0.41 ug/L # 63
35) ethyl tert-butyl ether	8.594	59	2770	0.38 ug/L 92
37) 1,1-dichloroethane	8.153	63	1154	0.32 ug/L 91
38) chloroprene	8.258	53	903	0.33 ug/L 76
39) acrylonitrile	7.571	53	1255	1.88 ug/L 89
42) 2,2-dichloropropane	8.882	77	1441	0.43 ug/L 85
43) cis-1,2-dichloroethene	8.882	96	1007	0.48 ug/L # 37
46) bromochloromethane	9.212	128	323	0.33 ug/L # 44
48) chloroform	9.260	83	1250	0.37 ug/L 91
49) Tert-Butyl Formate	9.291	59	768	0.41 ug/L # 82
54) 1,1,1-trichloroethane	9.511	97	1198	0.38 ug/L 89
57) tert-amyl methyl ether	9.983	73	3140	0.46 ug/L 87
62) carbon tetrachloride	9.705	117	806	0.34 ug/L 91
63) 1,1-dichloropropene	9.679	75	536	0.26 ug/L # 52
64) hexane	7.886	57	709	0.37 ug/L # 47
65) benzene	9.952	78	2575	0.40 ug/L 93
68) 1,2-dichloroethane	9.962	62	663	0.33 ug/L # 42
69) trichloroethene	10.675	95	623	0.44 ug/L 80
75) 1,2-dichloropropane	10.937	63	663	0.40 ug/L 90
76) methylcyclohexane	10.890	83	886	0.29 ug/L 78
77) dibromomethane	11.095	93	290	0.31 ug/L # 71
78) bromodichloromethane	11.226	83	837	0.43 ug/L 89
79) cis-1,3-dichloropropene	11.687	75	808	0.37 ug/L 69
82) toluene	12.044	92	1220	0.36 ug/L 98
84) trans-1,3-dichloropropene	12.264	75	795	0.41 ug/L 83
86) 1,1,2-trichloroethane	12.453	83	432	0.41 ug/L 82
89) tetrachloroethene	12.636	166	697	0.45 ug/L 68
90) 1,3-dichloropropane	12.647	76	796	0.40 ug/L 90

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112236.d
 Acq On : 29 Jul 2015 8:42 pm
 Operator : ximenac
 Sample : IC4810-0.5
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Aug 06 13:48:29 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

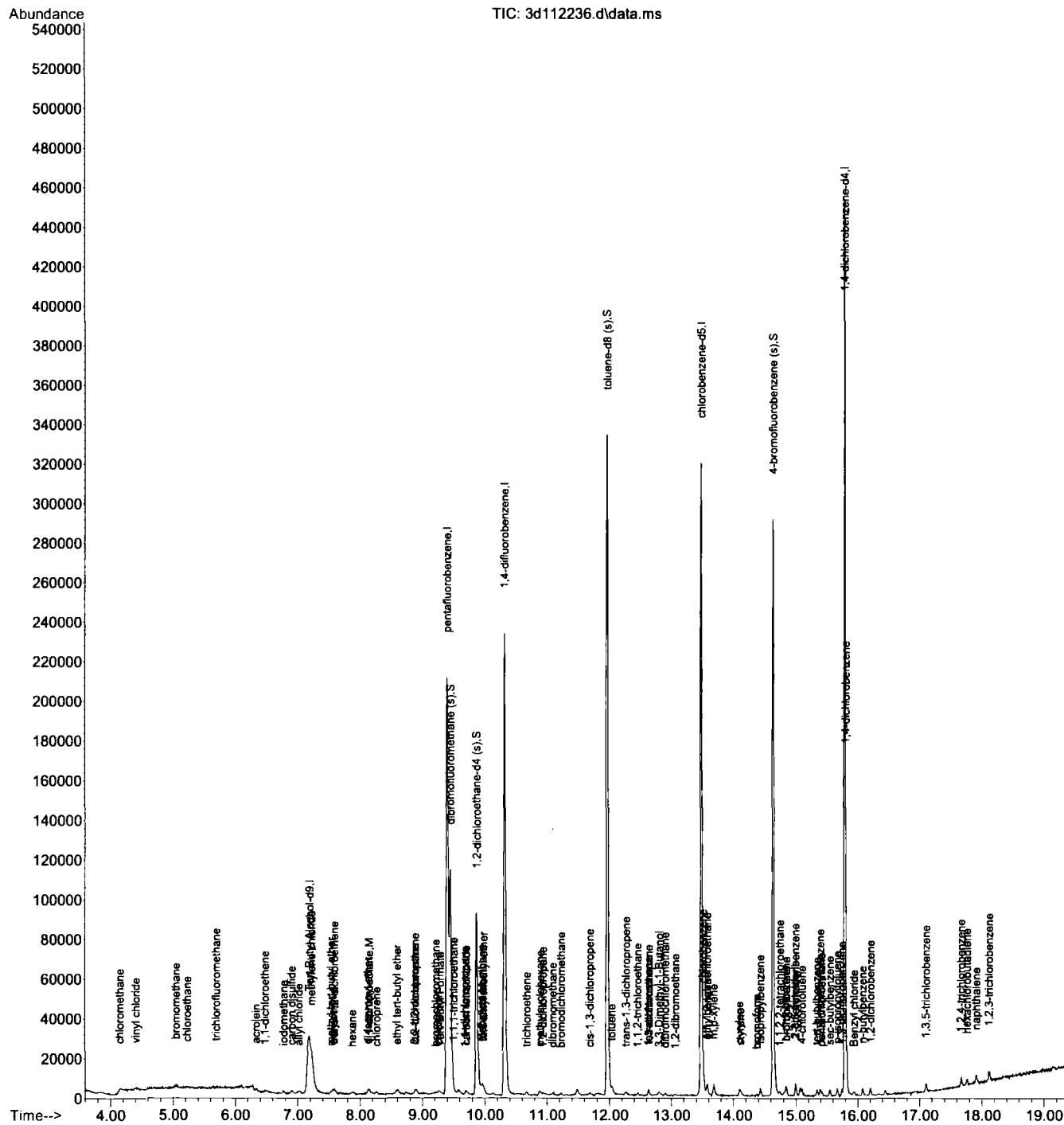
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
92) 3,3-Dimethyl-1-Butanol	12.814	57	950	5.30	ug/L #	90
93) dibromochloromethane	12.904	129	503	0.33	ug/L #	59
94) 1,2-dibromoethane	13.066	107	489	0.37	ug/L	89
96) chlorobenzene	13.512	112	1533	0.40	ug/L	92
97) 1,1,1,2-tetrachloroethane	13.569	131	539	0.34	ug/L	91
98) ethylbenzene	13.580	91	2807	0.43	ug/L	94
99) m,p-xylene	13.680	106	1837	0.75	ug/L	88
100) o-xylene	14.094	106	949	0.39	ug/L	98
101) styrene	14.120	104	1435	0.37	ug/L	90
103) bromoform	14.372	173	449	0.40	ug/L	91
105) isopropylbenzene	14.435	105	2697	0.38	ug/L	97
107) bromobenzene	14.843	156	848	0.43	ug/L #	53
109) 1,1,2,2-tetrachloroethane	14.744	83	859	0.42	ug/L	72
112) n-propylbenzene	14.849	91	2878	0.36	ug/L	94
114) 2-chlorotoluene	14.996	126	727	0.42	ug/L #	58
115) 4-chlorotoluene	15.095	126	691	0.41	ug/L #	73
116) 1,3,5-trimethylbenzene	14.990	105	2331	0.39	ug/L	84
117) tert-butylbenzene	15.342	119	1703	0.34	ug/L	83
118) pentachloroethane	15.420	167	406	0.31	ug/L	86
119) 1,2,4-trimethylbenzene	15.394	105	2258	0.37	ug/L	92
120) sec-butylbenzene	15.557	105	2714	0.34	ug/L	96
121) 1,3-dichlorobenzene	15.745	146	1527	0.39	ug/L	90
122) p-isopropyltoluene	15.677	119	2660	0.39	ug/L	97
123) 1,4-dichlorobenzene	15.813	146	1649	0.41	ug/L	88
124) 1,2-dichlorobenzene	16.207	146	1561	0.37	ug/L	91
126) n-butylbenzene	16.081	92	1226	0.35	ug/L	92
129) 1,3,5-trichlorobenzene	17.108	180	1680	0.39	ug/L	90
130) 1,2,4-trichlorobenzene	17.669	180	1563	0.35	ug/L	96
131) hexachlorobutadiene	17.764	225	687	0.33	ug/L	94
132) naphthalene	17.911	128	3428	0.35	ug/L	92
133) 1,2,3-trichlorobenzene	18.115	180	1803	0.39	ug/L	83
135) Benzyl chloride	15.939	91	1862	0.49	ug/L #	86

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112236.d
Acq On : 29 Jul 2015 8:42 pm
Operator : ximenac
Sample : IC4810-0.5
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 21 Sample Multiplier: 1

Quant Time: Aug 06 13:48:29 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4801-4810\
 Data File : 3d112237.d
 Acq On : 29 Jul 2015 9:10 pm
 Operator : ximenac
 Sample : IC4810-1
 Misc : MS88759, V3D4810, 5, , , 1
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Aug 06 13:49:00 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.173	65	97101	500.00	ug/L	0.00
4) pentafluorobenzene	9.396	168	208847	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	236603	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	197557	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	125183	50.00	ug/L	0.00

System Monitoring Compounds

50) dibromofluoromethane (s)	9.443	113	80599	44.62	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 89.24%	
51) 1,2-dichloroethane-d4 (s)	9.868	65	74529	39.96	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 79.92%	
80) toluene-d8 (s)	11.965	98	261711	51.94	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 103.88%	
106) 4-bromofluorobenzene (s)	14.634	95	98509	46.50	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 93.00%	

Target Compounds

					Qvalue	
2) 1,4-dioxane	11.058	88	321	19.81	ug/L	# 39
9) dichlorodifluoromethane	3.833	85	3408m	0.80	ug/L	
10) chloromethane	4.137	50	2804	0.90	ug/L	95
11) vinyl chloride	4.394	62	2869	0.84	ug/L	91
12) bromomethane	5.039	94	2094	1.25	ug/L	98
13) chloroethane	5.207	64	1340	1.06	ug/L	87
16) trichlorofluoromethane	5.663	101	2816	0.83	ug/L	79
18) ethyl ether	6.067	74	948	1.14	ug/L	83
21) acrolein	6.324	56	311518	768.98	ug/L	100
22) 1,1-dichloroethene	6.491	61	3006	0.82	ug/L	91
24) allyl chloride	7.021	76	844	0.80	ug/L	# 41
25) acetonitrile	7.016	40	1710	7.46	ug/L	# 1
27) iodomethane	6.764	142	4119	0.95	ug/L	91
29) carbon disulfide	6.895	76	7493	0.84	ug/L	95
30) methylene chloride	7.204	84	2161	0.88	ug/L	96
31) methyl acetate	7.016	43	812	0.54	ug/L	# 63
32) methyl tert butyl ether	7.535	73	6136	0.90	ug/L	96
33) trans-1,2-dichloroethene	7.582	61	2738	0.86	ug/L	90
34) di-isopropyl ether	8.122	45	6515	0.84	ug/L	# 54
35) ethyl tert-butyl ether	8.583	59	5834	0.78	ug/L	89
37) 1,1-dichloroethane	8.153	63	3160	0.84	ug/L	99
38) chloroprene	8.263	53	2329	0.83	ug/L	88
39) acrylonitrile	7.566	53	2935	4.27	ug/L	94
42) 2,2-dichloropropane	8.882	77	3355	0.96	ug/L	93
43) cis-1,2-dichloroethene	8.887	96	2226	1.03	ug/L	87
44) propionitrile	8.982	54	1844	7.24	ug/L	69
46) bromochloromethane	9.202	128	994	0.98	ug/L	90
48) chloroform	9.249	83	2901	0.84	ug/L	97
49) Tert-Butyl Formate	9.270	59	1443	0.74	ug/L	# 41
52) freon 113	6.455	151	1653	0.93	ug/L	# 68
54) 1,1,1-trichloroethane	9.501	97	3005	0.93	ug/L	94
57) tert-amyl methyl ether	9.983	73	5744	0.81	ug/L	92
59) epichlorohydrin	11.598	57	741	5.57	ug/L	52
60) n-butyl alcohol	10.497	56	2075	54.71	ug/L	# 54
61) cyclohexane	9.569	84	3211	1.11	ug/L	82
62) carbon tetrachloride	9.705	117	2651	1.08	ug/L	95
63) 1,1-dichloropropene	9.684	75	1872	0.86	ug/L	88
64) hexane	7.870	57	1802	0.90	ug/L	75
65) benzene	9.952	78	6049	0.91	ug/L	98
66) heptane	10.125	57	919	0.86	ug/L	# 75
67) isopropyl acetate	9.878	43	2086	0.67	ug/L	# 1

Manual Integrations
APPROVED
(compounds with "m" flag)
Jessica Reitan-Chu
08/06/15 14:29

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112237.d
 Acq On : 29 Jul 2015 9:10 pm
 Operator : ximenac
 Sample : IC4810-1
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Aug 06 13:49:00 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

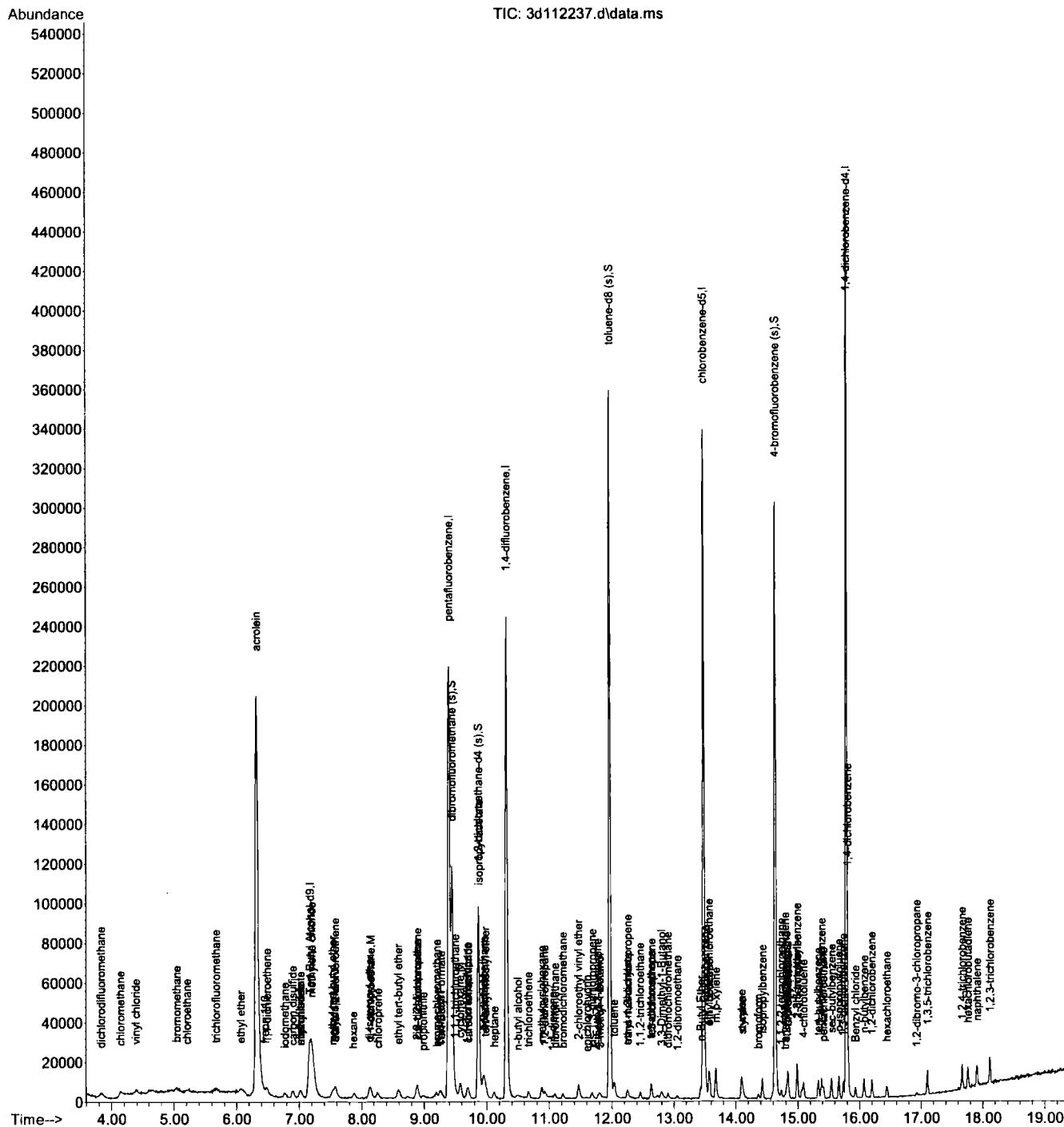
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
68) 1,2-dichloroethane	9.962	62	1745	0.83	ug/L	90
69) trichloroethene	10.670	95	1515	1.03	ug/L	93
73) 2-chloroethyl vinyl ether	11.472	63	4223	4.78	ug/L	95
75) 1,2-dichloropropane	10.927	63	1610	0.94	ug/L	89
76) methylcyclohexane	10.880	83	2926	0.93	ug/L	91
77) dibromomethane	11.089	93	905	0.93	ug/L	95
78) bromodichloromethane	11.221	83	1973	0.97	ug/L	95
79) cis-1,3-dichloropropene	11.687	75	2169	0.95	ug/L	75
81) 4-methyl-2-pentanone	11.776	58	437	0.78	ug/L #	10
82) toluene	12.038	92	3455	0.98	ug/L	86
83) 3-methyl-1-butanol	11.813	70	907	24.60	ug/L #	82
84) trans-1,3-dichloropropene	12.248	75	2028	0.99	ug/L	80
85) ethyl methacrylate	12.253	69	1658	1.05	ug/L	72
86) 1,1,2-trichloroethane	12.458	83	1211	1.11	ug/L	92
89) tetrachloroethene	12.631	166	1889	1.17	ug/L	85
90) 1,3-dichloropropane	12.641	76	1957	0.94	ug/L	98
92) 3,3-Dimethyl-1-Butanol	12.799	57	2325	12.35	ug/L #	91
93) dibromochloromethane	12.898	129	1506	0.95	ug/L	93
94) 1,2-dibromoethane	13.056	107	1190	0.86	ug/L	91
95) n-Butyl Ether	13.433	57	6600	1.01	ug/L #	93
96) chlorobenzene	13.512	112	3921	0.97	ug/L	99
97) 1,1,1,2-tetrachloroethane	13.564	131	1648	1.00	ug/L	77
98) ethylbenzene	13.569	91	6901	1.00	ug/L	94
99) m,p-xylene	13.674	106	5180	2.02	ug/L	97
100) o-xylene	14.089	106	2566	1.00	ug/L	100
101) styrene	14.109	104	4200	1.03	ug/L	92
103) bromoform	14.361	173	1106	0.94	ug/L	94
105) isopropylbenzene	14.429	105	7128	0.98	ug/L	98
107) bromobenzene	14.833	156	2029	1.03	ug/L #	69
109) 1,1,2,2-tetrachloroethane	14.733	83	2036	0.98	ug/L	86
110) trans-1,4-dichloro-2-b...	14.786	53	383	0.80	ug/L	92
111) 1,2,3-trichloropropene	14.807	110	431	0.95	ug/L	74
112) n-propylbenzene	14.838	91	7929	0.98	ug/L	99
114) 2-chlorotoluene	14.985	126	1626	0.94	ug/L	100
115) 4-chlorotoluene	15.090	126	1782	1.03	ug/L	92
116) 1,3,5-trimethylbenzene	14.985	105	5966	0.98	ug/L	95
117) tert-butylbenzene	15.331	119	4765	0.93	ug/L	90
118) pentachloroethane	15.420	167	1276	0.97	ug/L	94
119) 1,2,4-trimethylbenzene	15.384	105	6079	0.99	ug/L	96
120) sec-butylbenzene	15.551	105	8134	1.01	ug/L	98
121) 1,3-dichlorobenzene	15.735	146	3820	0.97	ug/L	93
122) p-isopropyltoluene	15.667	119	6957	1.00	ug/L	93
123) 1,4-dichlorobenzene	15.813	146	4380	1.07	ug/L	92
124) 1,2-dichlorobenzene	16.196	146	4185	0.99	ug/L	98
126) n-butylbenzene	16.076	92	3470	0.97	ug/L	81
128) 1,2-dibromo-3-chloropr...	16.925	157	520	0.94	ug/L	86
129) 1,3,5-trichlorobenzene	17.098	180	4284	0.99	ug/L	98
130) 1,2,4-trichlorobenzene	17.669	180	4286	0.96	ug/L	91
131) hexachlorobutadiene	17.759	225	2249	1.07	ug/L	95
132) naphthalene	17.905	128	8661	0.88	ug/L	99
133) 1,2,3-trichlorobenzene	18.115	180	4570	0.99	ug/L	93
134) hexachloroethane	16.448	119	1151	0.97	ug/L	95
135) Benzyl chloride	15.934	91	4283	1.12	ug/L	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112237.d
 Acq On : 29 Jul 2015 9:10 pm
 Operator : ximenac
 Sample : IC4810-1
 Misc : MS88759, V3D4810, 5, , , 1
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Aug 06 13:49:00 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration



Manual Integration Approval Summary

Page 1 of 1

Sample Number: V3D4810-IC4810 **Method:** SW846 8260C
Lab FileID: 3D112237.D **Analyst approved:** 08/06/15 13:52 Henny Salim
Injection Time: 07/29/15 21:10 **Supervisor approved:** 08/06/15 14:29 Jessica Reitan-Chu

Parameter	CAS	Sig#	R.T. (min.)	Reason
Dichlorodifluoromethane	75-71-8		3.83	Poor instrument integration

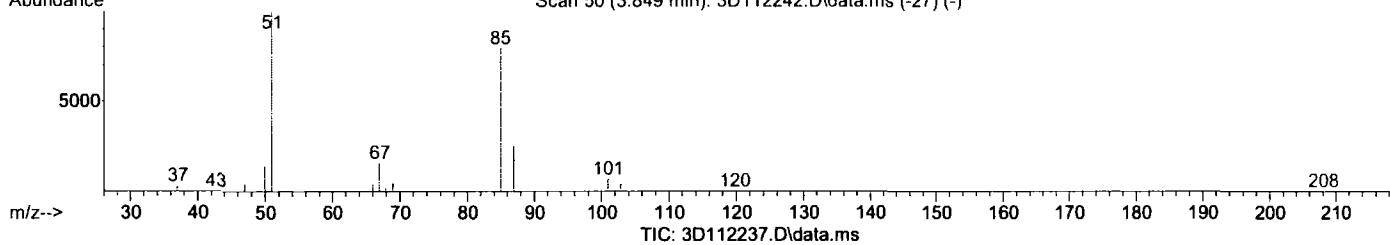
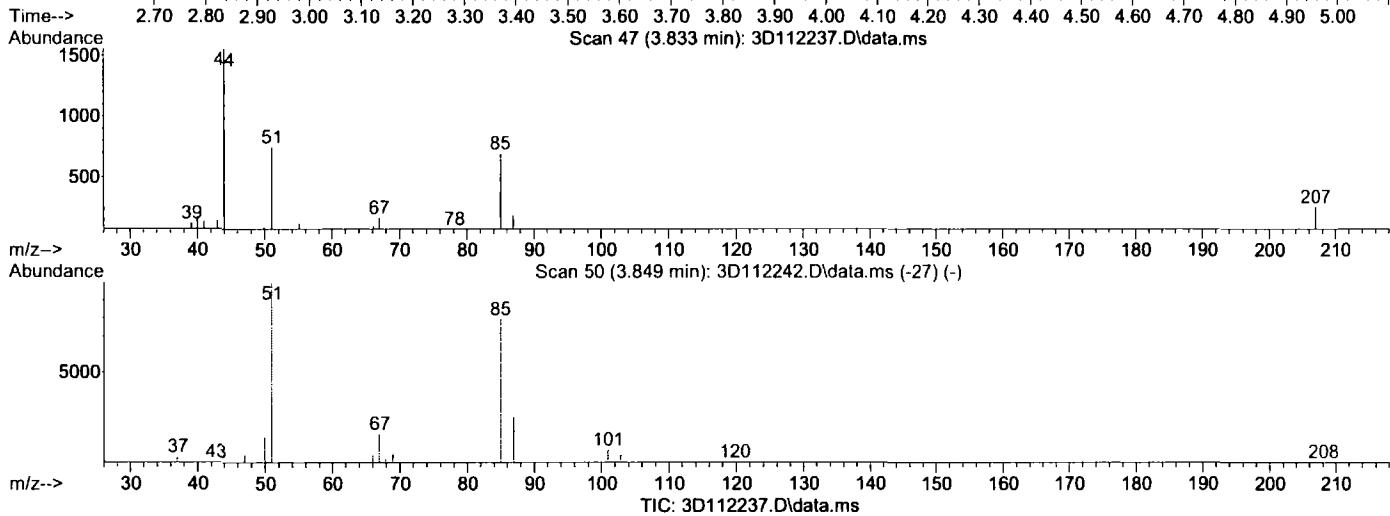
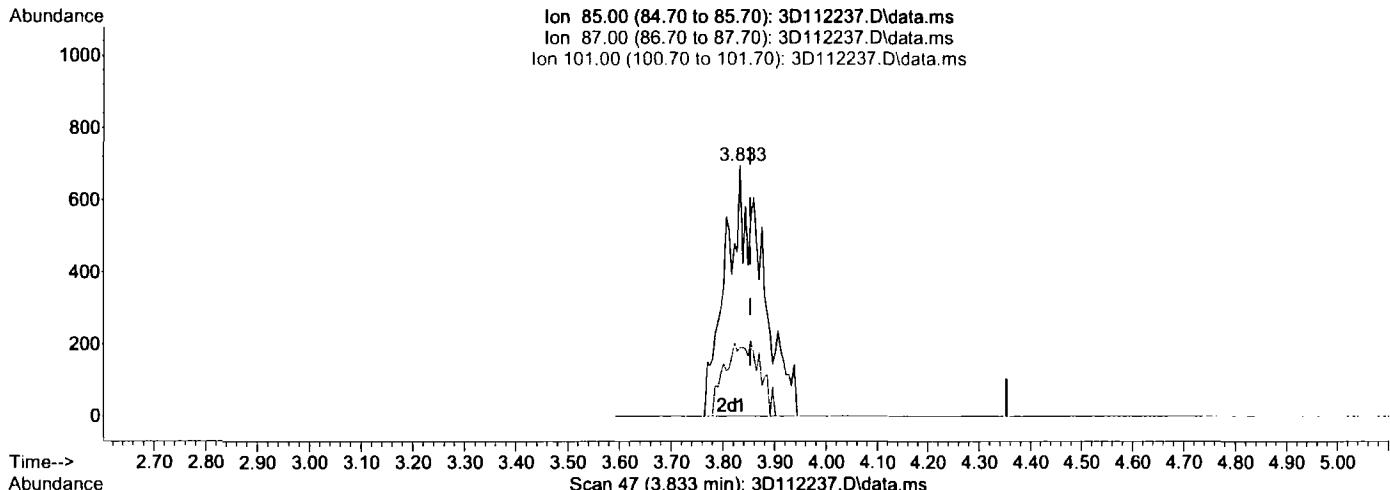
7631

7

Quantitation Report (Qedit)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112237.D
 Acq On : 29 Jul 2015 9:10 pm
 Operator : ximenac
 Sample : IC4810-1
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Jul 30 08:37:55 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration



(9) dichlorodifluoromethane

3.833min (-0.019) 0.80ug/L m

response 3408

Ion	Exp%	Act%
85.00	100	100
87.00	31.30	27.30
101.00	8.10	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Manual Integrations	
APPROVED	
(compounds with "m" flag)	
Jessica Reitan-Chu	08/06/15 14:29

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112238.d
Acq On : 29 Jul 2015 9:37 pm
Operator : ximenac
Sample : IC4810-2
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 23 Sample Multiplier: 1

Quant Time: Aug 06 13:49:15 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.173	65	91896	500.00	ug/L	0.00
4) pentafluorobenzene	9.396	168	197423	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	221121	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	180387	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.793	152	114767	50.00	ug/L	0.00

System Monitoring Compounds

50) dibromofluoromethane (s)	9.443	113	74452	43.60	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	87.20%
51) 1,2-dichloroethane-d4 (s)	9.863	65	70397	39.93	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	79.86%
80) toluene-d8 (s)	11.965	98	244288	51.88	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	103.76%
106) 4-bromofluorobenzene (s)	14.634	95	87810	45.21	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.42%

Target Compounds

					Value
2) 1,4-dioxane	11.048	88	694	45.25	ug/L # 60
3) tertiary butyl alcohol	7.283	59	2143m	10.37	ug/L
8) chlorodifluoromethane	3.844	51	5524	1.41	ug/L 93
9) dichlorodifluoromethane	3.844	85	7278	1.82	ug/L 96
10) chloromethane	4.148	50	5016	1.70	ug/L 98
11) vinyl chloride	4.394	62	5371	1.67	ug/L 97
12) bromomethane	5.034	94	3528	2.23	ug/L 81
13) chloroethane	5.202	64	2287	1.91	ug/L 95
16) trichlorofluoromethane	5.674	101	5823	1.81	ug/L 85
18) ethyl ether	6.072	74	1425	1.81	ug/L 83
21) acrolein	6.318	56	595837	1555.92	ug/L 97
22) 1,1-dichloroethene	6.486	61	4751	1.38	ug/L 97
24) allyl chloride	7.011	76	1423	1.43	ug/L # 85
25) acetonitrile	7.000	40	2912	13.44	ug/L # 36
27) iodomethane	6.769	142	6893	1.68	ug/L 97
28) iso-butyl alcohol	9.705	41	926	13.21	ug/L # 72
29) carbon disulfide	6.890	76	12180	1.45	ug/L 96
30) methylene chloride	7.205	84	3549	1.52	ug/L 97
31) methyl acetate	7.016	43	1975	1.38	ug/L # 88
32) methyl tert butyl ether	7.540	73	10043	1.56	ug/L 96
33) trans-1,2-dichloroethene	7.582	61	4499	1.49	ug/L 92
34) di-isopropyl ether	8.127	45	10502	1.44	ug/L # 56
35) ethyl tert-butyl ether	8.594	59	10092	1.42	ug/L 96
37) 1,1-dichloroethane	8.143	63	5031	1.42	ug/L 93
38) chloroprene	8.253	53	4066	1.53	ug/L 89
39) acrylonitrile	7.556	53	4607	7.09	ug/L 89
41) ethyl acetate	8.888	45	479	2.22	ug/L # 39
42) 2,2-dichloropropane	8.888	77	5181	1.57	ug/L 90
43) cis-1,2-dichloroethene	8.888	96	3646	1.78	ug/L # 81
44) propionitrile	8.966	54	3462	14.39	ug/L 72
45) methyl acrylate	8.977	55	1966	1.24	ug/L 73
46) bromochloromethane	9.197	128	1530	1.59	ug/L # 81
47) tetrahydrofuran	9.244	42	1230	1.88	ug/L # 61
48) chloroform	9.249	83	5020	1.55	ug/L 94
49) Tert-Butyl Formate	9.276	59	2632	1.44	ug/L # 88
52) freon 113	6.444	151	2895	1.72	ug/L 79
53) methacrylonitrile	9.139	41	1630	1.67	ug/L 62
54) 1,1,1-trichloroethane	9.491	97	4812	1.57	ug/L 97
57) tert-amyl methyl ether	9.978	73	10083	1.51	ug/L 97
59) epichlorohydrin	11.593	57	1322	10.63	ug/L 68
60) n-butyl alcohol	10.471	56	2972	83.84	ug/L 90

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4810-4810\
 Data File : 3d112238.d
 Acq On : 29 Jul 2015 9:37 pm
 Operator : ximenac
 Sample : IC4810-2
 Misc : MS88759, V3D4810.5,,,1
 ALS Vial : 23 Sample Multiplier: 1

Quant Time: Aug 06 13:49:15 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
61) cyclohexane	9.574	84	5021	1.86	ug/L	94
62) carbon tetrachloride	9.711	117	4206	1.83	ug/L	96
63) 1,1-dichloropropene	9.679	75	3058	1.51	ug/L	84
64) hexane	7.876	57	3186	1.69	ug/L	71
65) benzene	9.941	78	10496	1.69	ug/L	97
66) heptane	10.125	57	1727	1.73	ug/L	82
67) isopropyl acetate	9.879	43	4629	1.60	ug/L	62
68) 1,2-dichloroethane	9.962	62	3007	1.52	ug/L	89
69) trichloroethene	10.665	95	2636	1.91	ug/L	94
73) 2-chloroethyl vinyl ether	11.462	63	7676	9.29	ug/L	94
74) methyl methacrylate	10.948	41	2823	1.26	ug/L	89
75) 1,2-dichloropropane	10.922	63	2567	1.60	ug/L	94
76) methylcyclohexane	10.880	83	5176	1.76	ug/L	93
77) dibromomethane	11.084	93	1592	1.76	ug/L	93
78) bromodichloromethane	11.215	83	3129	1.64	ug/L	92
79) cis-1,3-dichloropropene	11.682	75	3854	1.80	ug/L	94
81) 4-methyl-2-pentanone	11.776	58	920	1.75	ug/L	96
82) toluene	12.039	92	5701	1.73	ug/L	95
83) 3-methyl-1-butanol	11.813	70	1458	42.31	ug/L	84
84) trans-1,3-dichloropropene	12.248	75	3201	1.68	ug/L	87
85) ethyl methacrylate	12.243	69	2635	1.79	ug/L	83
86) 1,1,2-trichloroethane	12.458	83	1827	1.80	ug/L	92
87) 2-hexanone	12.663	58	632m	1.51	ug/L	
89) tetrachloroethene	12.631	166	2996	2.03	ug/L	99
90) 1,3-dichloropropane	12.636	76	3564	1.87	ug/L	96
91) butyl acetate	12.725	56	1247	1.53	ug/L	# 56
92) 3,3-Dimethyl-1-Butanol	12.799	57	3466	20.16	ug/L	# 91
93) dibromochloromethane	12.898	129	2576	1.78	ug/L	91
94) 1,2-dibromoethane	13.051	107	2138	1.68	ug/L	94
95) n-Butyl Ether	13.428	57	10714	1.79	ug/L	# 89
96) chlorobenzene	13.512	112	6372	1.73	ug/L	88
97) 1,1,1,2-tetrachloroethane	13.570	131	2712	1.80	ug/L	94
98) ethylbenzene	13.570	91	10810	1.72	ug/L	99
99) m,p-xylene	13.674	106	8457	3.61	ug/L	99
100) o-xylene	14.089	106	4328	1.84	ug/L	88
101) styrene	14.104	104	6752	1.81	ug/L	93
103) bromoform	14.361	173	1953	1.82	ug/L	83
105) isopropylbenzene	14.429	105	11312	1.70	ug/L	98
107) bromobenzene	14.833	156	3361	1.86	ug/L	89
109) 1,1,2,2-tetrachloroethane	14.734	83	3182	1.67	ug/L	95
110) trans-1,4-dichloro-2-b...	14.786	53	628	1.44	ug/L	# 80
111) 1,2,3-trichloropropane	14.807	110	646	1.56	ug/L	79
112) n-propylbenzene	14.838	91	13012	1.75	ug/L	99
114) 2-chlorotoluene	14.985	126	2828	1.77	ug/L	# 75
115) 4-chlorotoluene	15.090	126	2646	1.67	ug/L	88
116) 1,3,5-trimethylbenzene	14.990	105	9596	1.71	ug/L	93
117) tert-butylbenzene	15.331	119	8045	1.71	ug/L	96
118) pentachloroethane	15.415	167	2010	1.67	ug/L	90
119) 1,2,4-trimethylbenzene	15.384	105	9434	1.68	ug/L	93
120) sec-butylbenzene	15.546	105	13151	1.78	ug/L	99
121) 1,3-dichlorobenzene	15.735	146	6184	1.71	ug/L	94
122) p-isopropyltoluene	15.667	119	11115	1.75	ug/L	99
123) 1,4-dichlorobenzene	15.814	146	6495	1.74	ug/L	97
124) 1,2-dichlorobenzene	16.196	146	6835	1.77	ug/L	95
126) n-butylbenzene	16.076	92	5866	1.78	ug/L	92
128) 1,2-dibromo-3-chloropr...	16.925	157	883	1.74	ug/L	95
129) 1,3,5-trichlorobenzene	17.098	180	7345	1.85	ug/L	98
130) 1,2,4-trichlorobenzene	17.664	180	7027	1.71	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112238.d
Acq On : 29 Jul 2015 9:37 pm
Operator : ximenac
Sample : IC4810-2
Misc : MS88759,V3D4810,5,,,,1
ALS Vial : 23 Sample Multiplier: 1

Quant Time: Aug 06 13:49:15 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration

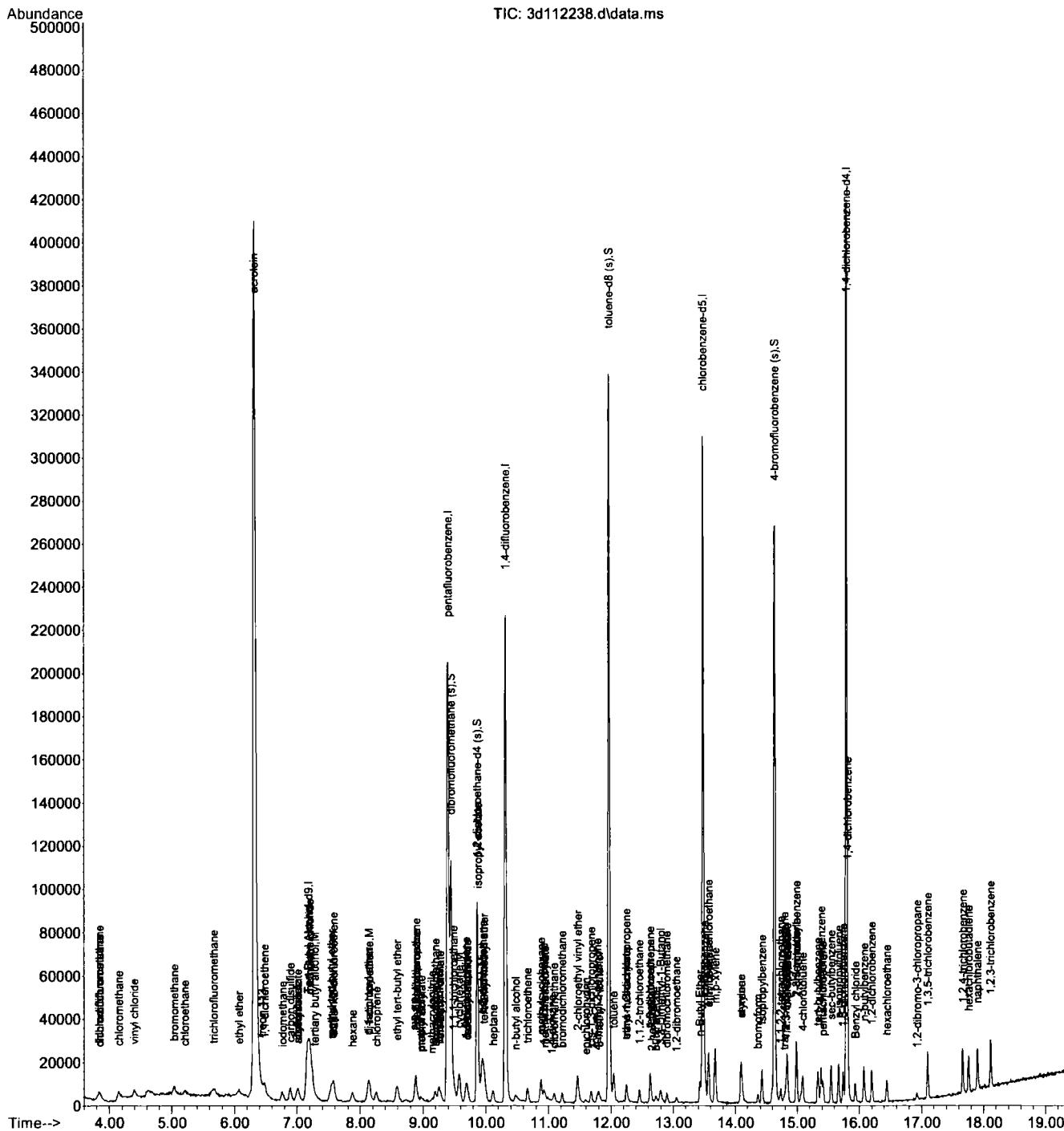
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
131) hexachlorobutadiene	17.759	225	3791	1.96	ug/L	94
132) naphthalene	17.900	128	14293	1.58	ug/L	98
133) 1,2,3-trichlorobenzene	18.110	180	7641	1.80	ug/L	91
134) hexachloroethane	16.443	119	1902	1.76	ug/L	91
135) Benzyl chloride	15.929	91	7260	2.06	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112238.d
Acq On : 29 Jul 2015 9:37 pm
Operator : ximenac
Sample : IC4810-2
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 23 Sample Multiplier: 1

Quant Time: Aug 06 13:49:15 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration



7.6.4

Manual Integration Approval Summary

Page 1 of 1

Sample Number: V3D4810-IC4810 **Method:** SW846 8260C
Lab FileID: 3D112238.D **Analyst approved:** 08/06/15 13:52 Henny Salim
Injection Time: 07/29/15 21:37 **Supervisor approved:** 08/06/15 14:29 Jessica Reitan-Chu

Parameter	CAS	Sig#	R.T. (min.)	Reason
Tert Butyl Alcohol	75-65-0		7.28	Poor instrument integration
2-Hexanone	591-78-6		12.66	Poor instrument integration

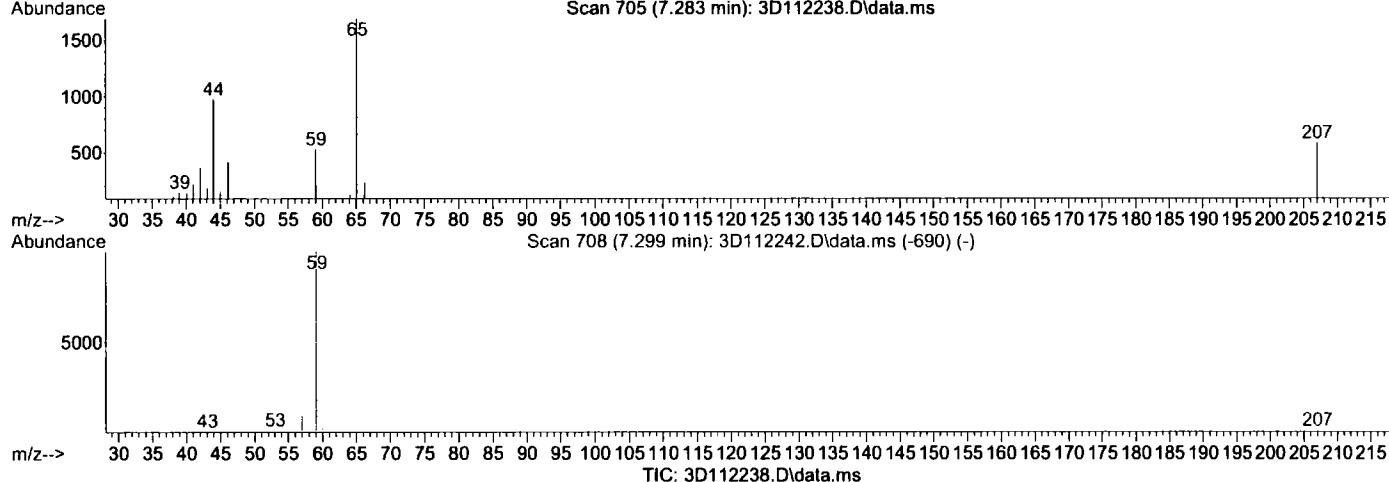
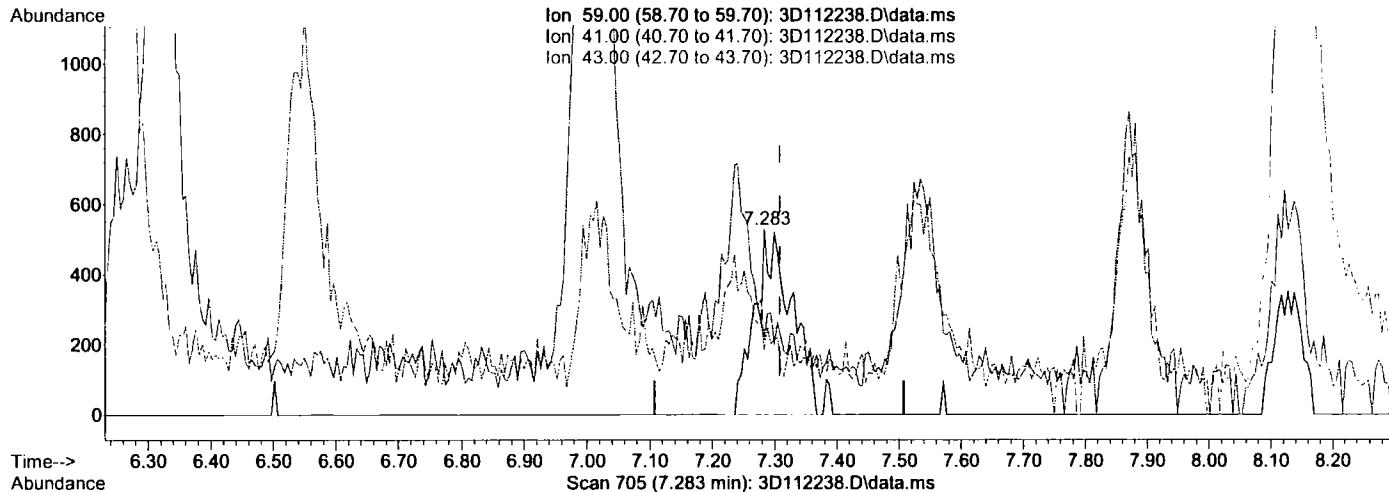
14.61

L

Quantitation Report (Qedit)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112238.D
 Acq On : 29 Jul 2015 9:37 pm
 Operator : ximenac
 Sample : IC4810-2
 Misc : MS88759,V3D4810.5,,,1
 ALS Vial : 23 Sample Multiplier: 1

Quant Time: Jul 30 08:37:57 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration



(3) tertiary butyl alcohol (M)

7.283min (-0.024) 10.37ug/L m

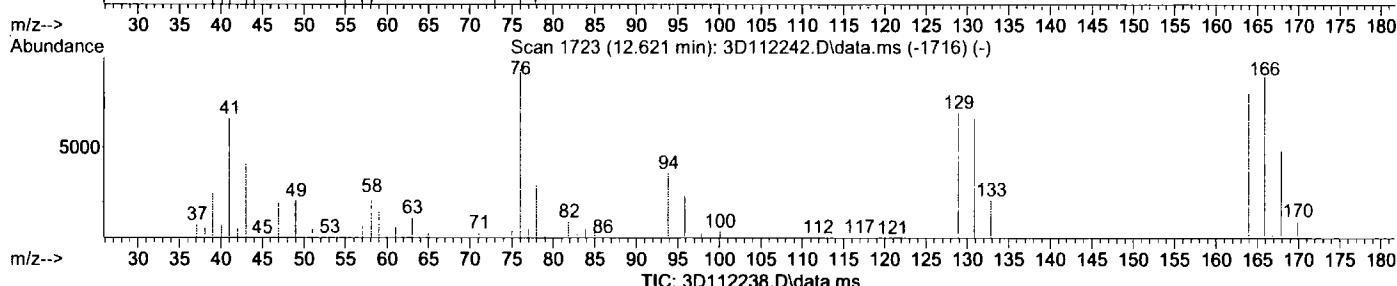
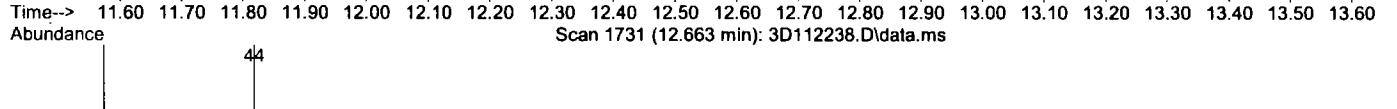
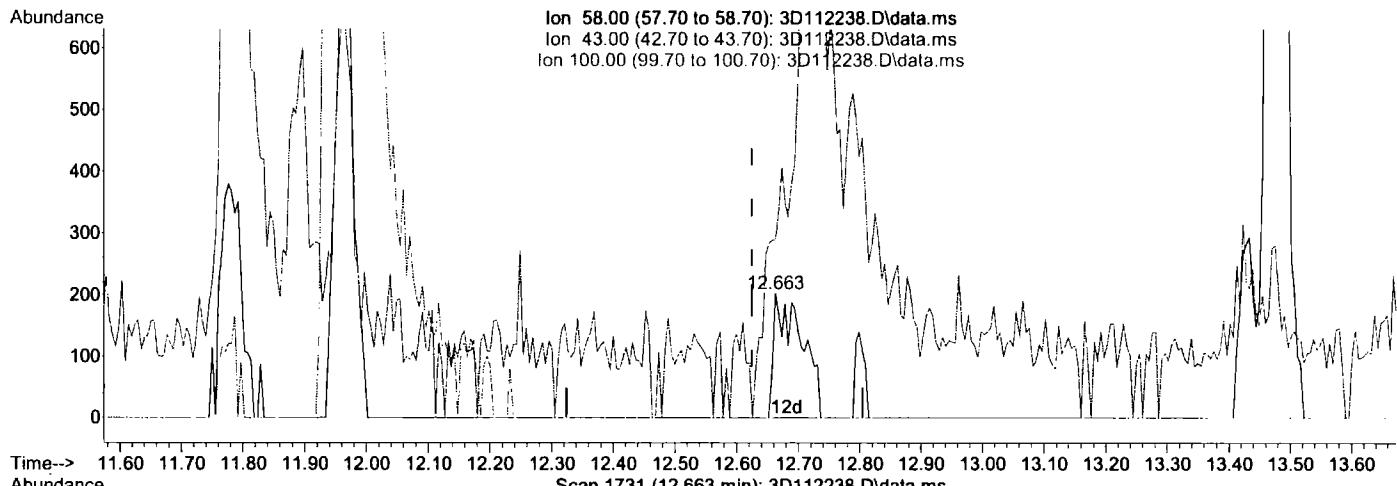
response 2143

Ion	Exp%	Act%
59.00	100	100
41.00	30.20	41.02
43.00	16.60	34.59
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112238.D
 Acq On : 29 Jul 2015 9:37 pm
 Operator : ximenac
 Sample : IC4810-2
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 23 Sample Multiplier: 1

Quant Time: Jul 30 08:37:57 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration



(87) 2-hexanone

12.663min (+0.039) 1.51ug/L m

response 632

Ion	Exp%	Act%
58.00	100	100
43.00	183.00	142.57#
100.00	14.90	0.00
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4810-4810\
 Data File : 3d112239.d
 Acq On : 29 Jul 2015 10:04 pm
 Operator : ximenac
 Sample : IC4810-5
 Misc : MS88759, V3D4810, 5, , , 1
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Jul 30 11:45:40 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.173	65	103357	500.00	ug/L	0.00
4) pentafluorobenzene	9.396	168	208424	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	236372	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	193775	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	122426	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev (Min)
50) dibromofluoromethane (s)	9.443	113	79004	43.82	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery =	87.64%		
51) 1,2-dichloroethane-d4 (s)	9.863	65	75160	40.38	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery =	80.76%		
80) toluene-d8 (s)	11.965	98	257202	51.10	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery =	102.20%		
106) 4-bromofluorobenzene (s)	14.629	95	95829	46.26	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	92.52%		

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-dioxane	11.042	88	1992	115.48	ug/L	# 65
3) tertiary butyl alcohol	7.294	59	5490	23.62	ug/L	85
8) chlorodifluoromethane	3.854	51	14241	3.45	ug/L	89
9) dichlorodifluoromethane	3.838	85	20494	4.85	ug/L	98
10) chloromethane	4.148	50	14329	4.61	ug/L	95
11) vinyl chloride	4.405	62	15435	4.55	ug/L	98
12) bromomethane	5.044	94	9842	5.89	ug/L	91
13) chloroethane	5.212	64	6641	5.24	ug/L	96
16) trichlorofluoromethane	5.663	101	16831	4.95	ug/L	96
18) ethyl ether	6.082	74	3733	4.49	ug/L	84
21) acrolein	6.329	56	16708	41.33	ug/L	94
22) 1,1-dichloroethene	6.491	61	12228	3.36	ug/L	94
23) acetone	6.554	43	2497	4.20	ug/L	92
24) allyl chloride	7.016	76	3802	3.61	ug/L	# 64
25) acetonitrile	6.989	40	7985	34.91	ug/L	87
27) iodomethane	6.769	142	17419	4.03	ug/L	96
28) iso-butyl alcohol	9.700	41	2677	36.16	ug/L	92
29) carbon disulfide	6.895	76	30912	3.48	ug/L	97
30) methylene chloride	7.204	84	8855	3.60	ug/L	97
31) methyl acetate	7.010	43	5821	3.86	ug/L	# 93
32) methyl tert butyl ether	7.529	73	26062	3.82	ug/L	95
33) trans-1,2-dichloroethene	7.582	61	10922	3.42	ug/L	93
34) di-isopropyl ether	8.127	45	28139	3.65	ug/L	# 66
35) ethyl tert-butyl ether	8.589	59	28689	3.83	ug/L	97
36) 2-butanone	8.877	72	551	3.21	ug/L	# 40
37) 1,1-dichloroethane	8.143	63	13258	3.54	ug/L	97
38) chloroprene	8.253	53	10153	3.61	ug/L	88
39) acrylonitrile	7.545	53	13036	19.01	ug/L	95
40) vinyl acetate	8.153	86	820	3.41	ug/L	# 1
41) ethyl acetate	8.882	45	1135	4.99	ug/L	# 1
42) 2,2-dichloropropane	8.882	77	13031	3.74	ug/L	97
43) cis-1,2-dichloroethene	8.882	96	8516	3.94	ug/L	89
44) propionitrile	8.956	54	9368	36.88	ug/L	# 57
45) methyl acrylate	8.982	55	5927	3.55	ug/L	95
46) bromochloromethane	9.191	128	4132	4.07	ug/L	# 80
47) tetrahydrofuran	9.244	42	2497	3.62	ug/L	94
48) chloroform	9.249	83	12212	3.56	ug/L	94
49) Tert-Butyl Formate	9.281	59	7337	3.79	ug/L	# 86
52) freon 113	6.455	151	7988	4.50	ug/L	79
53) methacrylonitrile	9.144	41	3379	3.28	ug/L	86
54) 1,1,1-trichloroethane	9.490	97	12500	3.87	ug/L	87

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4810-4810\
 Data File : 3d112239.d
 Acq On : 29 Jul 2015 10:04 pm
 Operator : ximenac
 Sample : IC4810-5
 Misc : MS88759, V3D4810, 5, , , 1
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Jul 30 11:45:40 2015
 Quant Method : C:\msdchem\1\METHODS\3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
57) tert-amyl methyl ether	9.978	73	26879	3.81	ug/L	97
59) epichlorohydrin	11.588	57	3344	25.15	ug/L	93
60) n-butyl alcohol	10.455	56	9220	243.33	ug/L	96
61) cyclohexane	9.574	84	12866	4.45	ug/L	87
62) carbon tetrachloride	9.711	117	11265	4.60	ug/L	92
63) 1,1-dichloropropene	9.684	75	8394	3.87	ug/L	92
64) hexane	7.875	57	7999	3.98	ug/L	91
65) benzene	9.941	78	26303	3.96	ug/L	96
66) heptane	10.114	57	4552	4.27	ug/L	93
67) isopropyl acetate	9.873	43	11841	3.83	ug/L	80
68) 1,2-dichloroethane	9.957	62	7828	3.71	ug/L	97
69) trichloroethene	10.665	95	6095	4.13	ug/L	92
72) 2-nitropropane	11.436	41	2599	4.46	ug/L	94
73) 2-chloroethyl vinyl ether	11.457	63	21267	24.08	ug/L	97
74) methyl methacrylate	10.943	41	8395	3.49	ug/L	90
75) 1,2-dichloropropane	10.927	63	6904	4.03	ug/L	98
76) methylcyclohexane	10.880	83	13545	4.31	ug/L	94
77) dibromomethane	11.084	93	4071	4.21	ug/L	85
78) bromodichloromethane	11.215	83	8288	4.07	ug/L	92
79) cis-1,3-dichloropropene	11.677	75	9561	4.19	ug/L	96
81) 4-methyl-2-pentanone	11.771	58	2559	4.56	ug/L	95
82) toluene	12.038	92	14664	4.17	ug/L	91
83) 3-methyl-1-butanol	11.797	70	3909	106.12	ug/L	93
84) trans-1,3-dichloropropene	12.243	75	8786	4.30	ug/L	92
85) ethyl methacrylate	12.243	69	7405	4.70	ug/L	94
86) 1,1,2-trichloroethane	12.453	83	4712	4.33	ug/L	95
87) 2-hexanone	12.641	58	2039	4.54	ug/L	91
89) tetrachloroethene	12.631	166	7517	4.74	ug/L	99
90) 1,3-dichloropropane	12.636	76	8725	4.25	ug/L	97
91) butyl acetate	12.710	56	3621	4.13	ug/L	86
92) 3,3-Dimethyl-1-Butanol	12.793	57	8670	46.95	ug/L #	95
93) dibromochloromethane	12.898	129	6864	4.41	ug/L	94
94) 1,2-dibromoethane	13.050	107	5896	4.32	ug/L	99
95) n-Butyl Ether	13.428	57	27368	4.25	ug/L	95
96) chlorobenzene	13.507	112	16872	4.26	ug/L	93
97) 1,1,2,2-tetrachloroethane	13.564	131	6983	4.31	ug/L	95
98) ethylbenzene	13.564	91	27948	4.15	ug/L	100
99) m,p-xylene	13.669	106	22094	8.77	ug/L	93
100) o-xylene	14.089	106	11324	4.48	ug/L	92
101) styrene	14.099	104	17957	4.47	ug/L	96
103) bromoform	14.361	173	5321	4.62	ug/L	97
105) isopropylbenzene	14.424	105	29522	4.17	ug/L	99
107) bromobenzene	14.828	156	8206	4.25	ug/L	92
108) cyclohexanone	14.592	55	9785	51.82	ug/L	96
109) 1,1,2,2-tetrachloroethane	14.733	83	8583	4.22	ug/L	98
110) trans-1,4-dichloro-2-b...	14.775	53	2042	4.38	ug/L	81
111) 1,2,3-trichloropropane	14.807	110	1984	4.48	ug/L	97
112) n-propylbenzene	14.838	91	33004	4.17	ug/L	100
114) 2-chlorotoluene	14.985	126	7464	4.39	ug/L	90
115) 4-chlorotoluene	15.085	126	7495	4.44	ug/L	91
116) 1,3,5-trimethylbenzene	14.985	105	25327	4.24	ug/L	95
117) tert-butylbenzene	15.331	119	19703	3.92	ug/L	100
118) pentachloroethane	15.415	167	5254	4.10	ug/L	97
119) 1,2,4-trimethylbenzene	15.378	105	25772	4.30	ug/L	99
120) sec-butylbenzene	15.546	105	33972	4.32	ug/L	99
121) 1,3-dichlorobenzene	15.735	146	16173	4.20	ug/L	98
122) p-isopropyltoluene	15.667	119	29093	4.29	ug/L	98
123) 1,4-dichlorobenzene	15.813	146	17166	4.30	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4810-4810\
 Data File : 3d112239.d
 Acq On : 29 Jul 2015 10:04 pm
 Operator : ximenac
 Sample : IC4810-5
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Jul 30 11:45:40 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

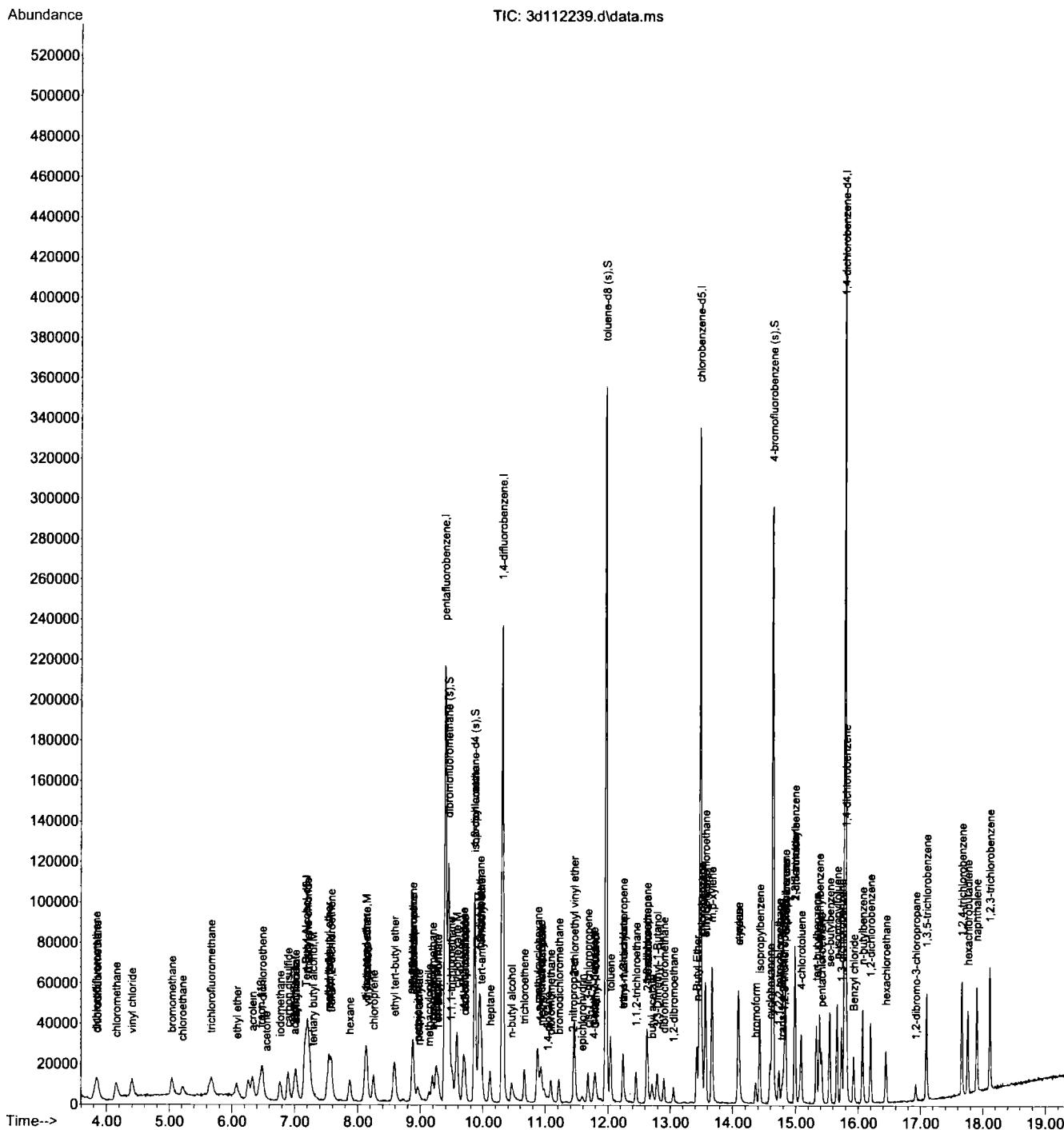
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.196	146	17639	4.27	ug/L	98
126) n-butylbenzene	16.070	92	14766	4.20	ug/L	97
128) 1,2-dibromo-3-chloropr...	16.930	157	2382	4.39	ug/L	82
129) 1,3,5-trichlorobenzene	17.098	180	18081	4.26	ug/L	99
130) 1,2,4-trichlorobenzene	17.664	180	18349	4.19	ug/L	98
131) hexachlorobutadiene	17.759	225	9315	4.52	ug/L	92
132) naphthalene	17.900	128	38147	3.95	ug/L	98
133) 1,2,3-trichlorobenzene	18.110	180	19180	4.23	ug/L	95
134) hexachloroethane	16.443	119	5108	4.42	ug/L	98
135) Benzyl chloride	15.929	91	17026	4.54	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112239.d
Acq On : 29 Jul 2015 10:04 pm
Operator : ximenac
Sample : IC4810-5
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 24 Sample Multiplier: 1

Quant Time: Jul 30 11:45:40 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration



M3D4810.M Thu Aug 06 13:50:21 2015 3D

Page : 4

Manual Integrations
APPROVED
(compounds with "Pm" flag)
Jessica Reitan-Chu
08/06/15 14:29

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112240.d
 Acq On : 29 Jul 2015 10:31 pm
 Operator : ximenac
 Sample : IC4810-10
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 30 11:48:54 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.168	65	94749	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	201206	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	224714	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	185472	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	118281	50.00	ug/L	0.00

System Monitoring Compounds

50) dibromofluoromethane (s)	9.443	113	75974	43.66	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 87.32%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	70388	39.18	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 78.36%	
80) toluene-d8 (s)	11.960	98	246270	51.46	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 102.92%	
106) 4-bromofluorobenzene (s)	14.629	95	91078	45.50	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 91.00%	

Target Compounds

					Qvalue
2) 1,4-dioxane	11.037	88	3736	236.26	ug/L
3) tertiary butyl alcohol	7.299	59	9830	46.14	ug/L
8) chlorodifluoromethane	3.849	51	24666	6.19	ug/L
9) dichlorodifluoromethane	3.838	85	38339	9.40	ug/L
10) chloromethane	4.148	50	25990	8.66	ug/L
11) vinyl chloride	4.405	62	29135	8.89	ug/L
12) bromomethane	5.050	94	17972	11.14	ug/L
13) chloroethane	5.217	64	12105	9.90	ug/L
16) trichlorofluoromethane	5.673	101	31073	9.47	ug/L
18) ethyl ether	6.072	74	7336	9.13	ug/L
21) acrolein	6.324	56	29882	76.56	ug/L
22) 1,1-dichloroethene	6.491	61	23852	6.79	ug/L
23) acetone	6.549	43	4844	8.43	ug/L
24) allyl chloride	7.016	76	7520	7.39	ug/L #
25) acetonitrile	6.989	40	13588	61.54	ug/L #
27) iodomethane	6.769	142	34492	8.27	ug/L
28) iso-butyl alcohol	9.700	41	4644	64.99	ug/L
29) carbon disulfide	6.895	76	59937	6.99	ug/L
30) methylene chloride	7.210	84	17711	7.47	ug/L
31) methyl acetate	7.005	43	9525	6.53	ug/L
32) methyl tert butyl ether	7.535	73	52934	8.05	ug/L
33) trans-1,2-dichloroethene	7.577	61	21380	6.93	ug/L
34) di-isopropyl ether	8.127	45	49458	6.64	ug/L
35) ethyl tert-butyl ether	8.589	59	48644	6.72	ug/L
36) 2-butanone	8.872	72	1397m	8.42	ug/L
37) 1,1-dichloroethane	8.143	63	26481	7.32	ug/L
38) chloroprene	8.253	53	17145	6.32	ug/L
39) acrylonitrile	7.540	53	27491	41.52	ug/L
40) vinyl acetate	8.143	86	1532	6.60	ug/L #
41) ethyl acetate	8.877	45	1777	8.09	ug/L #
42) 2,2-dichloropropane	8.882	77	25506	7.59	ug/L
43) cis-1,2-dichloroethene	8.882	96	16793	8.04	ug/L
44) propionitrile	8.945	54	19260	78.54	ug/L
45) methyl acrylate	8.956	55	13110	8.13	ug/L
46) bromochloromethane	9.192	128	8533	8.71	ug/L
47) tetrahydrofuran	9.233	42	5013	7.53	ug/L
48) chloroform	9.244	83	23924	7.23	ug/L
49) Tert-Butyl Formate	9.275	59	12780	6.84	ug/L #
52) freon 113	6.460	151	13439	7.84	ug/L
53) methacrylonitrile	9.134	41	6903	6.94	ug/L
54) 1,1,1-trichloroethane	9.501	97	24746	7.94	ug/L

7.9.6

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112240.d
 Acq On : 29 Jul 2015 10:31 pm
 Operator : ximenac
 Sample : IC4810-10
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 30 11:48:54 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
57) tert-amyl methyl ether	9.973	73	45484	6.68	ug/L	97
59) epichlorohydrin	11.582	57	5766	45.62	ug/L	94
60) n-butyl alcohol	10.450	56	17409	483.28	ug/L	94
61) cyclohexane	9.569	84	25733	9.37	ug/L	82
62) carbon tetrachloride	9.705	117	21620	9.28	ug/L	99
63) 1,1-dichloropropene	9.679	75	16227	7.88	ug/L	92
64) hexane	7.876	57	13055	6.83	ug/L	95
65) benzene	9.941	78	51889	8.23	ug/L	97
66) heptane	10.114	57	7653	7.55	ug/L	92
67) isopropyl acetate	9.863	43	20448	6.95	ug/L	87
68) 1,2-dichloroethane	9.957	62	16233	8.10	ug/L	93
69) trichloroethene	10.660	95	12453	8.89	ug/L	98
72) 2-nitropropane	11.436	41	4289	7.74	ug/L	76
73) 2-chloroethyl vinyl ether	11.457	63	37588	44.77	ug/L	97
74) methyl methacrylate	10.927	41	17733	7.76	ug/L	97
75) 1,2-dichloropropane	10.917	63	13357	8.19	ug/L	99
76) methylcyclohexane	10.875	83	22551	7.55	ug/L	93
77) dibromomethane	11.084	93	8388	9.12	ug/L	96
78) bromodichloromethane	11.215	83	16668	8.61	ug/L	96
79) cis-1,3-dichloropropene	11.677	75	20153	9.28	ug/L	94
81) 4-methyl-2-pentanone	11.771	58	5014	9.39	ug/L	96
82) toluene	12.039	92	29423	8.80	ug/L	99
83) 3-methyl-1-butanol	11.792	70	7349	209.85	ug/L	91
84) trans-1,3-dichloropropene	12.238	75	18095	9.32	ug/L	91
85) ethyl methacrylate	12.227	69	15393	10.29	ug/L	98
86) 1,1,2-trichloroethane	12.447	83	9602	9.29	ug/L	92
87) 2-hexanone	12.636	58	3944	9.24	ug/L	# 79
89) tetrachloroethene	12.626	166	15228	10.03	ug/L	96
90) 1,3-dichloropropane	12.631	76	18003	9.17	ug/L	96
91) butyl acetate	12.704	56	6815	8.12	ug/L	# 75
92) 3,3-Dimethyl-1-Butanol	12.794	57	14747	83.43	ug/L	95
93) dibromochloromethane	12.898	129	14143	9.49	ug/L	95
94) 1,2-dibromoethane	13.050	107	12264	9.40	ug/L	94
95) n-Butyl Ether	13.423	57	53856	8.74	ug/L	99
96) chlorobenzene	13.507	112	33030	8.72	ug/L	95
97) 1,1,1,2-tetrachloroethane	13.569	131	14160	9.13	ug/L	98
98) ethylbenzene	13.564	91	55245	8.57	ug/L	98
99) m,p-xylene	13.669	106	43436	18.02	ug/L	99
100) o-xylene	14.089	106	21777	9.01	ug/L	88
101) styrene	14.099	104	36701	9.54	ug/L	96
103) bromoform	14.361	173	10922	9.90	ug/L	99
105) isopropylbenzene	14.424	105	58523	8.56	ug/L	97
107) bromobenzene	14.828	156	17244	9.25	ug/L	88
108) cyclohexanone	14.587	55	21965	120.40	ug/L	94
109) 1,1,2,2-tetrachloroethane	14.728	83	17033	8.67	ug/L	98
110) trans-1,4-dichloro-2-b...	14.775	53	3673	8.16	ug/L	93
111) 1,2,3-trichloropropane	14.802	110	3974	9.29	ug/L	98
112) n-propylbenzene	14.833	91	65204	8.52	ug/L	100
114) 2-chlorotoluene	14.980	126	14684	8.94	ug/L	96
115) 4-chlorotoluene	15.085	126	15054	9.23	ug/L	93
116) 1,3,5-trimethylbenzene	14.985	105	49921	8.65	ug/L	96
117) tert-butylbenzene	15.331	119	39986	8.23	ug/L	98
118) pentachloroethane	15.415	167	10172	8.21	ug/L	95
119) 1,2,4-trimethylbenzene	15.378	105	50402	8.70	ug/L	99
120) sec-butylbenzene	15.546	105	66528	8.75	ug/L	98
121) 1,3-dichlorobenzene	15.735	146	33232	8.94	ug/L	98
122) p-isopropyltoluene	15.661	119	57504	8.78	ug/L	97
123) 1,4-dichlorobenzene	15.813	146	34937	9.06	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112240.d
 Acq On : 29 Jul 2015 10:31 pm
 Operator : ximenac
 Sample : IC4810-10
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 30 11:48:54 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

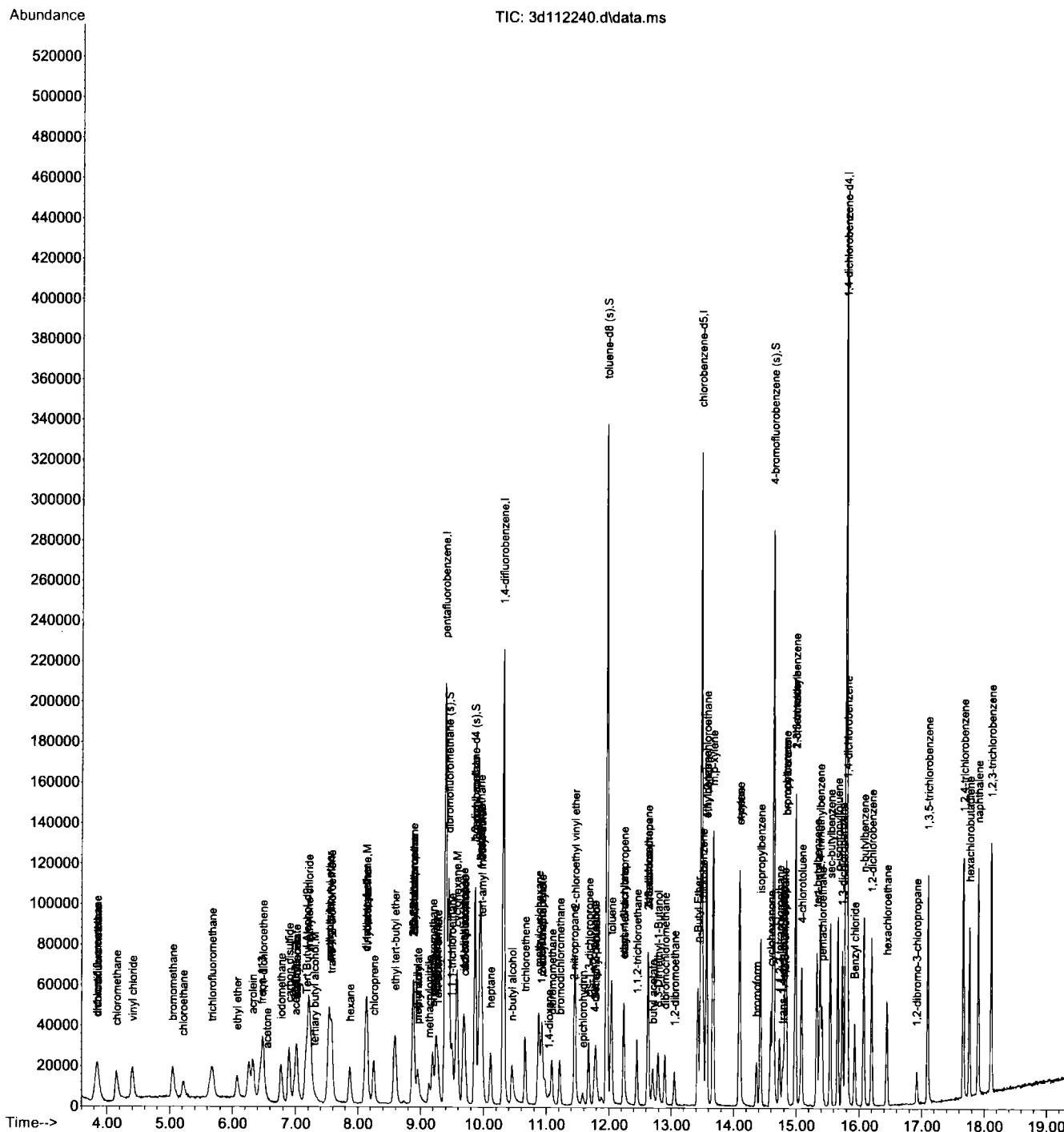
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
124) 1,2-dichlorobenzene	16.196	146	35274	8.84	ug/L	99
126) n-butylbenzene	16.070	92	30437	8.96	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.925	157	4754	9.07	ug/L	90
129) 1,3,5-trichlorobenzene	17.098	180	37404	9.12	ug/L	99
130) 1,2,4-trichlorobenzene	17.664	180	38620	9.12	ug/L	96
131) hexachlorobutadiene	17.759	225	18465	9.28	ug/L	98
132) naphthalene	17.900	128	79663	8.55	ug/L	99
133) 1,2,3-trichlorobenzene	18.110	180	38953	8.89	ug/L	97
134) hexachloroethane	16.443	119	10337	9.26	ug/L	95
135) Benzyl chloride	15.929	91	31103	8.58	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112240.d
Acq On : 29 Jul 2015 10:31 pm
Operator : ximenac
Sample : IC4810-10
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 30 11:48:54 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration



Manual Integration Approval Summary

Page 1 of 1

Sample Number: V3D4810-IC4810 **Method:** SW846 8260C
Lab FileID: 3D112240.D **Analyst approved:** 07/30/15 13:02 Henny Salim
Injection Time: 07/29/15 22:31 **Supervisor approved:** 08/06/15 14:29 Jessica Reitan-Chu

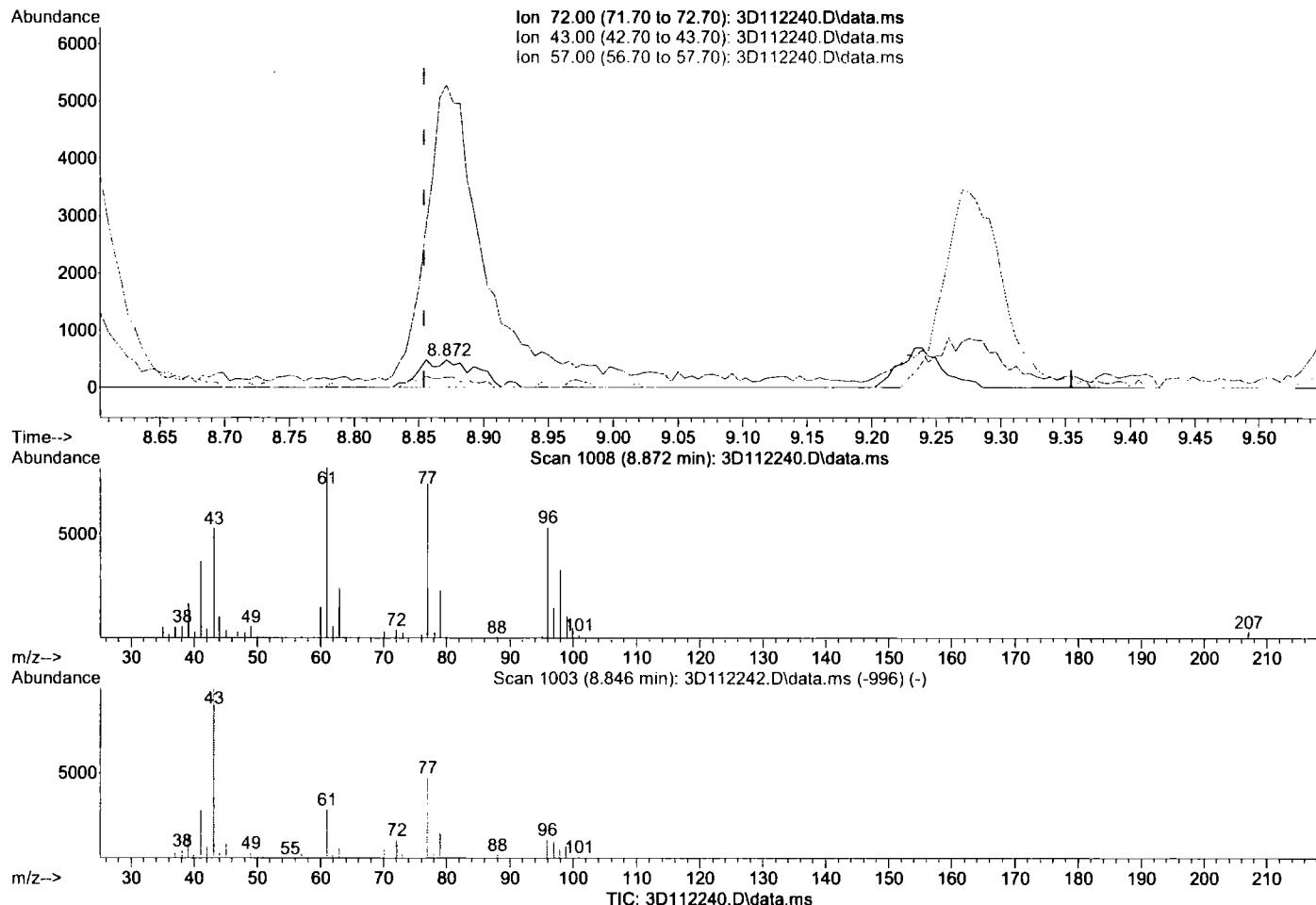
Parameter	CAS	Sig#	R.T. (min.)	Reason
2-Butanone (MEK)	78-93-3		8.87	Poor instrument integration

7661

Quantitation Report (Qedit)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112240.D
 Acq On : 29 Jul 2015 10:31 pm
 Operator : ximenac
 Sample : IC4810-10
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Jul 30 08:38:00 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration



(36) 2-butanone

8.872min (+0.018) 8.42ug/L m

response 1397

Ion	Exp%	Act%
72.00	100	100
43.00	1321.10	1109.66
57.00	32.30	37.65
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3D112241.D
 Acq On : 29 Jul 2015 10:58 pm
 Operator : ximenac
 Sample : IC4810-20
 Misc : MS88759,V3D4810.5,,,1
 ALS Vial : 26 Sample Multiplier: 1

Quant Time: Jul 30 11:50:38 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.168	65	96365	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	205871	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	233597	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	190298	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	121957	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
50) dibromofluoromethane (s)	9.443	113	78421	44.04	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery = 88.08%			
51) 1,2-dichloroethane-d4 (s)	9.863	65	72634	39.51	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery = 79.02%			
80) toluene-d8 (s)	11.960	98	257374	51.74	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery = 103.48%			
106) 4-bromofluorobenzene (s)	14.629	95	92946	45.04	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 90.08%			

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-dioxane	11.037	88	8448	525.29	ug/L	94
3) tertiary butyl alcohol	7.294	59	21383	98.68	ug/L	97
8) chlorodifluoromethane	3.844	51	56463	13.85	ug/L	96
9) dichlorodifluoromethane	3.844	85	77901	18.66	ug/L	99
10) chloromethane	4.153	50	53465	17.42	ug/L	98
11) vinyl chloride	4.405	62	58988	17.59	ug/L	95
12) bromomethane	5.044	94	37153	22.50	ug/L	97
13) chloroethane	5.212	64	23722	18.95	ug/L	95
16) trichlorofluoromethane	5.663	101	64098	19.08	ug/L	98
18) ethyl ether	6.072	74	15421	18.76	ug/L	97
21) acrolein	6.324	56	60721	152.06	ug/L	98
22) 1,1-dichloroethene	6.486	61	50149	13.94	ug/L	97
23) acetone	6.544	43	9878	16.80	ug/L	87
24) allyl chloride	7.011	76	16689	16.03	ug/L #	83
25) acetonitrile	6.974	40	30553	135.23	ug/L	88
27) iodomethane	6.764	142	74454	17.45	ug/L	97
28) iso-butyl alcohol	9.684	41	10148	138.79	ug/L	95
29) carbon disulfide	6.890	76	129241	14.72	ug/L	98
30) methylene chloride	7.205	84	37577	15.48	ug/L	99
31) methyl acetate	7.005	43	20757	13.92	ug/L	95
32) methyl tert butyl ether	7.530	73	112839	16.76	ug/L	97
33) trans-1,2-dichloroethene	7.577	61	46857	14.85	ug/L	92
34) di-isopropyl ether	8.122	45	110619	14.51	ug/L	90
35) ethyl tert-butyl ether	8.589	59	110696	14.94	ug/L	97
36) 2-butanone	8.856	72	3248	19.14	ug/L	67
37) 1,1-dichloroethane	8.143	63	57280	15.49	ug/L	97
38) chloroprene	8.248	53	38272	13.79	ug/L	93
39) acrylonitrile	7.535	53	59194	87.38	ug/L	99
40) vinyl acetate	8.138	86	4094	17.25	ug/L	52
41) ethyl acetate	8.872	45	4213	18.74	ug/L #	1
42) 2,2-dichloropropane	8.877	77	54222	15.77	ug/L	97
43) cis-1,2-dichloroethene	8.877	96	35402	16.57	ug/L	90
44) propionitrile	8.940	54	41971	167.28	ug/L	84
45) methyl acrylate	8.945	55	29812	18.07	ug/L	98
46) bromochloromethane	9.192	128	18677	18.63	ug/L #	82
47) tetrahydrofuran	9.228	42	10725	15.75	ug/L	97
48) chloroform	9.244	83	52546	15.52	ug/L	99
49) Tert-Butyl Formate	9.276	59	29933	15.66	ug/L	99
52) freon 113	6.460	151	29980	17.10	ug/L	81
53) methacrylonitrile	9.129	41	15856	15.59	ug/L	99
54) 1,1,1-trichloroethane	9.501	97	54834	17.20	ug/L	97

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3D112241.D
 Acq On : 29 Jul 2015 10:58 pm
 Operator : ximenac
 Sample : IC4810-20
 Misc : MS88759, V3D4810, 5, , , 1
 ALS Vial : 26 Sample Multiplier: 1

Quant Time: Jul 30 11:50:38 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
57) tert-amyl methyl ether	9.973	73	103703	14.88	ug/L	97
59) epichlorohydrin	11.577	57	12385	94.27	ug/L	93
60) n-butyl alcohol	10.445	56	37482	1000.95	ug/L	87
61) cyclohexane	9.574	84	53935	18.89	ug/L	87
62) carbon tetrachloride	9.705	117	48372	19.97	ug/L	98
63) 1,1-dichloropropene	9.674	75	36376	16.99	ug/L	94
64) hexane	7.870	57	28424	14.30	ug/L	94
65) benzene	9.936	78	112757	17.20	ug/L	98
66) heptane	10.109	57	16664	15.82	ug/L	96
67) isopropyl acetate	9.863	43	49101	16.06	ug/L	94
68) 1,2-dichloroethane	9.952	62	34905	16.75	ug/L	96
69) trichloroethene	10.660	95	27014	18.54	ug/L	91
72) 2-nitropropane	11.436	41	10648	18.49	ug/L	99
73) 2-chloroethyl vinyl ether	11.451	63	88704	101.63	ug/L	96
74) methyl methacrylate	10.927	41	41085	17.30	ug/L	91
75) 1,2-dichloropropane	10.922	63	29465	17.38	ug/L	98
76) methylcyclohexane	10.880	83	51606	16.62	ug/L	94
77) dibromomethane	11.084	93	18582	19.44	ug/L	92
78) bromodichloromethane	11.210	83	36643	18.21	ug/L	96
79) cis-1,3-dichloropropene	11.672	75	45351	20.09	ug/L	95
81) 4-methyl-2-pentanone	11.771	58	11182	20.15	ug/L	98
82) toluene	12.039	92	66864	19.23	ug/L	99
83) 3-methyl-1-butanol	11.787	70	14832	407.43	ug/L	91
84) trans-1,3-dichloropropene	12.233	75	40244	19.95	ug/L	99
85) ethyl methacrylate	12.227	69	34164	21.96	ug/L	98
86) 1,1,2-trichloroethane	12.448	83	21411	19.93	ug/L	97
87) 2-hexanone	12.626	58	9318	21.00	ug/L	98
89) tetrachloroethene	12.626	166	35050	22.49	ug/L	98
90) 1,3-dichloropropane	12.631	76	39330	19.52	ug/L	98
91) butyl acetate	12.699	56	15344	17.81	ug/L	96
92) 3,3-Dimethyl-1-Butanol	12.788	57	31726	174.95	ug/L	98
93) dibromochloromethane	12.898	129	31662	20.70	ug/L	96
94) 1,2-dibromoethane	13.045	107	27208	20.32	ug/L	96
95) n-Butyl Ether	13.423	57	118568	18.75	ug/L	99
96) chlorobenzene	13.507	112	74600	19.20	ug/L	97
97) 1,1,1,2-tetrachloroethane	13.564	131	32140	20.20	ug/L	97
98) ethylbenzene	13.559	91	122689	18.55	ug/L	99
99) m,p-xylene	13.669	106	95400	38.58	ug/L	100
100) o-xylene	14.083	106	49837	20.09	ug/L	90
101) styrene	14.099	104	80880	20.50	ug/L	93
103) bromoform	14.361	173	24626	21.75	ug/L	99
105) isopropylbenzene	14.424	105	130640	18.53	ug/L	98
107) bromobenzene	14.823	156	37524	19.51	ug/L	94
108) cyclohexanone	14.587	55	49760	264.53	ug/L	98
109) 1,1,2,2-tetrachloroethane	14.728	83	37265	18.39	ug/L	98
110) trans-1,4-dichloro-2-b...	14.775	53	8343	17.97	ug/L	95
111) 1,2,3-trichloropropane	14.802	110	8484	19.24	ug/L	90
112) n-propylbenzene	14.833	91	142828	18.11	ug/L	100
114) 2-chlorotoluene	14.980	126	32656	19.28	ug/L	92
115) 4-chlorotoluene	15.080	126	32386	19.27	ug/L	98
116) 1,3,5-trimethylbenzene	14.985	105	111035	18.65	ug/L	95
117) tert-butylbenzene	15.331	119	91815	18.32	ug/L	96
118) pentachloroethane	15.415	167	23555	18.43	ug/L	96
119) 1,2,4-trimethylbenzene	15.378	105	111069	18.59	ug/L	100
120) sec-butylbenzene	15.546	105	150463	19.19	ug/L	98
121) 1,3-dichlorobenzene	15.735	146	73542	19.18	ug/L	99
122) p-isopropyltoluene	15.662	119	129273	19.15	ug/L	98
123) 1,4-dichlorobenzene	15.814	146	75030	18.88	ug/L	97

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4810-4810\
 Data File : 3D112241.D
 Acq On : 29 Jul 2015 10:58 pm
 Operator : ximenac
 Sample : IC4810-20
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 26 Sample Multiplier: 1

Quant Time: Jul 30 11:50:38 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

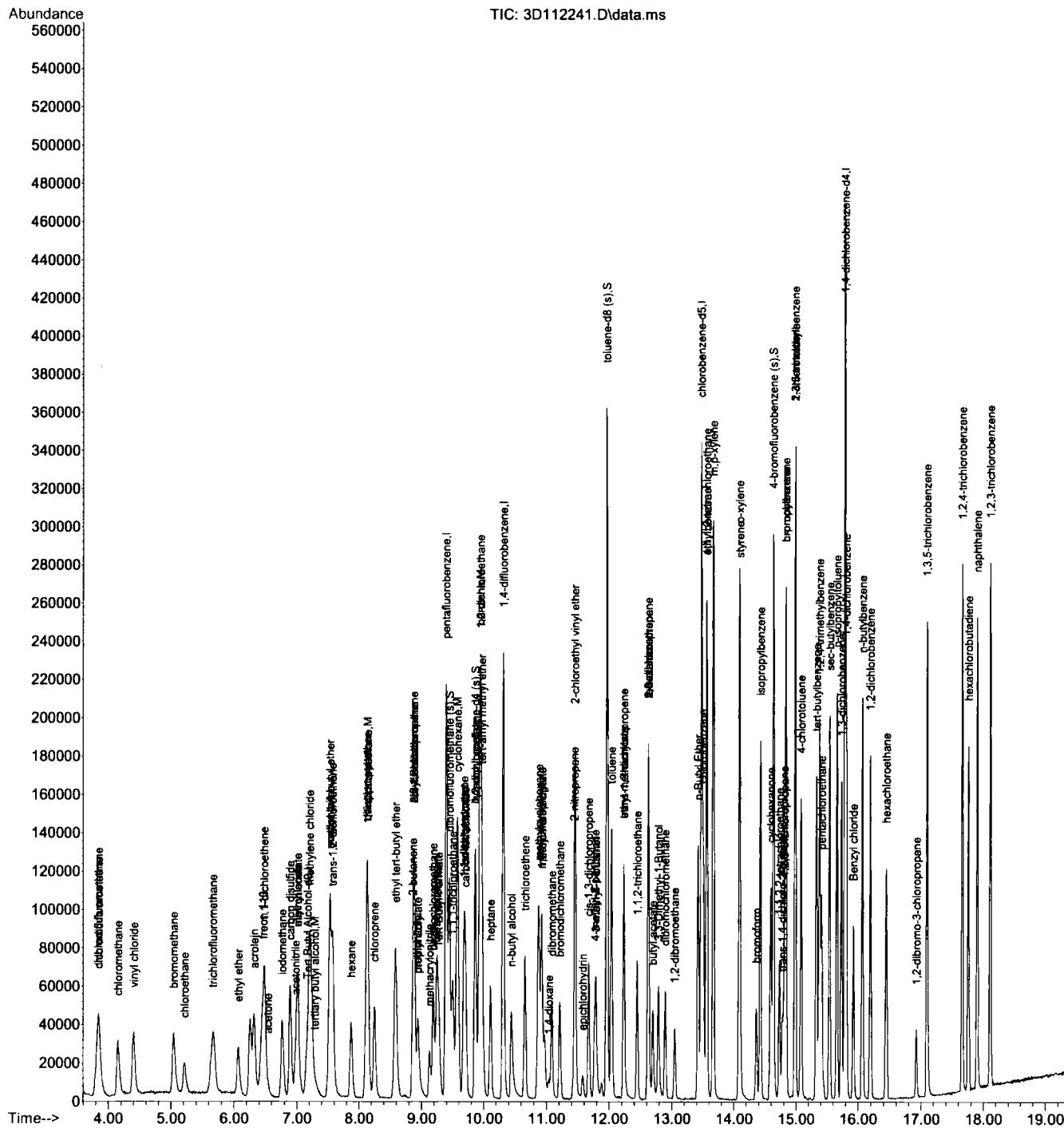
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.196	146	77329	18.79	ug/L	97
126) n-butylbenzene	16.065	92	67003	19.14	ug/L	98
128) 1,2-dibromo-3-chloropr...	16.925	157	11109	20.55	ug/L	98
129) 1,3,5-trichlorobenzene	17.098	180	84834	20.07	ug/L	97
130) 1,2,4-trichlorobenzene	17.659	180	85540	19.60	ug/L	98
131) hexachlorobutadiene	17.759	225	39564	19.29	ug/L	98
132) naphthalene	17.900	128	178644	18.59	ug/L	99
133) 1,2,3-trichlorobenzene	18.110	180	87582	19.38	ug/L	99
134) hexachloroethane	16.443	119	24382	21.19	ug/L	98
135) Benzyl chloride	15.924	91	66494	17.79	ug/L	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3D112241.D
Acq On : 29 Jul 2015 10:58 pm
Operator : ximenac
Sample : IC4810-20
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 26 Sample Multiplier: 1

Quant Time: Jul 30 11:50:38 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112242.d
 Acq On : 29 Jul 2015 11:25 pm
 Operator : ximenac
 Sample : ICC4810-50
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Jul 30 11:52:18 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.178	65	96354	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	213808	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	238297	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	193300	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	122120	50.00	ug/L	0.00

System Monitoring Compounds

50) dibromofluoromethane (s)	9.443	113	82616	44.67	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 89.34%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	74743	39.15	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 78.30%	
80) toluene-d8 (s)	11.960	98	262476	51.72	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 103.44%	
106) 4-bromofluorobenzene (s)	14.629	95	94748	45.85	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 91.70%	

Target Compounds

					Qvalue	
2) 1,4-dioxane	11.037	88	21435	1332.96	ug/L	88
3) tertiary butyl alcohol	7.299	59	51882	239.46	ug/L	96
8) chlorodifluoromethane	3.854	51	144662	34.17	ug/L	98
9) dichlorodifluoromethane	3.849	85	193205	44.57	ug/L	98
10) chloromethane	4.164	50	145170	45.53	ug/L	99
11) vinyl chloride	4.415	62	161354	46.33	ug/L	99
12) bromomethane	5.050	94	96095	56.03	ug/L	97
13) chloroethane	5.212	64	62712	48.25	ug/L	97
16) trichlorofluoromethane	5.674	101	168107	48.19	ug/L	99
18) ethyl ether	6.072	74	38059	44.58	ug/L	98
21) acrolein	6.324	56	145068	349.79	ug/L	98
22) 1,1-dichloroethene	6.486	61	132542	35.49	ug/L	97
23) acetone	6.544	43	20305	33.26	ug/L	83
24) allyl chloride	7.016	76	43027	39.79	ug/L #	81
25) acetonitrile	6.979	40	75584	322.12	ug/L	91
27) iodomethane	6.769	142	194884	43.98	ug/L	98
28) iso-butyl alcohol	9.685	41	25463	335.31	ug/L	100
29) carbon disulfide	6.895	76	334343	36.68	ug/L	99
30) methylene chloride	7.205	84	97033	38.49	ug/L	95
31) methyl acetate	7.000	43	53986	34.85	ug/L	98
32) methyl tert butyl ether	7.535	73	282066	40.35	ug/L	96
33) trans-1,2-dichloroethene	7.577	61	119916	36.59	ug/L	94
34) di-isopropyl ether	8.122	45	292146	36.90	ug/L	95
35) ethyl tert-butyl ether	8.584	59	294220	38.24	ug/L	98
36) 2-butanone	8.846	72	7359	41.76	ug/L	75
37) 1,1-dichloroethane	8.138	63	145923	37.98	ug/L	99
38) chloroprene	8.248	53	105722	36.68	ug/L	96
39) acrylonitrile	7.530	53	142501	202.55	ug/L	97
40) vinyl acetate	8.133	86	10726	43.51	ug/L	64
41) ethyl acetate	8.861	45	9145	39.18	ug/L	95
42) 2,2-dichloropropane	8.882	77	137502	38.51	ug/L	98
43) cis-1,2-dichloroethene	8.877	96	89025	40.13	ug/L	92
44) propionitrile	8.940	54	101157	388.20	ug/L	98
45) methyl acrylate	8.940	55	70423	41.10	ug/L	98
46) bromochloromethane	9.186	128	46329	44.49	ug/L #	83
47) tetrahydrofuran	9.228	42	24184	34.20	ug/L	97
48) chloroform	9.244	83	136324	38.76	ug/L	97
49) Tert-Butyl Formate	9.276	59	79581	40.10	ug/L #	89
52) freon 113	6.450	151	80037	43.95	ug/L	86
53) methacrylonitrile	9.124	41	38770	36.70	ug/L	93
54) 1,1,1-trichloroethane	9.496	97	140366	42.39	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112242.d
 Acq On : 29 Jul 2015 11:25 pm
 Operator : ximenac
 Sample : ICC4810-50
 Misc : MS88759,V3D4810,5,,,,1
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Jul 30 11:52:18 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
57) tert-amyl methyl ether	9.973	73	273411	37.77	ug/L	97
59) epichlorohydrin	11.577	57	29046	216.73	ug/L	95
60) n-butyl alcohol	10.440	56	95617	2503.08	ug/L	93
61) cyclohexane	9.569	84	141247	48.48	ug/L	84
62) carbon tetrachloride	9.706	117	126210	51.07	ug/L	99
63) 1,1-dichloropropene	9.674	75	91905	42.07	ug/L	93
64) hexane	7.876	57	76756	37.85	ug/L	95
65) benzene	9.936	78	284170	42.48	ug/L	98
66) heptane	10.109	57	43171	40.18	ug/L	97
67) isopropyl acetate	9.858	43	122747	39.34	ug/L	97
68) 1,2-dichloroethane	9.952	62	86550	40.70	ug/L	98
69) trichloroethene	10.655	95	68269	45.94	ug/L	95
72) 2-nitropropane	11.431	41	23129	39.37	ug/L	97
73) 2-chloroethyl vinyl ether	11.451	63	217123	243.86	ug/L	98
74) methyl methacrylate	10.922	41	98478	40.65	ug/L	96
75) 1,2-dichloropropane	10.917	63	74464	43.07	ug/L	99
76) methylcyclohexane	10.875	83	141516	44.67	ug/L	95
77) dibromomethane	11.084	93	44776	45.93	ug/L	89
78) bromodichloromethane	11.210	83	92134	44.88	ug/L	94
79) cis-1,3-dichloropropene	11.672	75	111103	48.24	ug/L	94
81) 4-methyl-2-pentanone	11.761	58	26308	46.47	ug/L	98
82) toluene	12.033	92	165473	46.65	ug/L	98
83) 3-methyl-1-butanol	11.787	70	36375	979.49	ug/L	88
84) trans-1,3-dichloropropene	12.233	75	99123	48.16	ug/L	91
85) ethyl methacrylate	12.227	69	83527	52.63	ug/L	95
86) 1,1,2-trichloroethane	12.448	83	51777	47.25	ug/L	95
87) 2-hexanone	12.621	58	20948	46.29	ug/L	98
89) tetrachloroethene	12.626	166	85256	53.86	ug/L	99
90) 1,3-dichloropropane	12.626	76	95746	46.78	ug/L	96
91) butyl acetate	12.694	56	39546	45.20	ug/L	93
92) 3,3-Dimethyl-1-Butanol	12.788	57	78103	423.99	ug/L	100
93) dibromochloromethane	12.893	129	78012	50.21	ug/L	96
94) 1,2-dibromoethane	13.045	107	64300	47.27	ug/L	100
95) n-Butyl Ether	13.423	57	292584	45.55	ug/L	98
96) chlorobenzene	13.507	112	181711	46.05	ug/L	96
97) 1,1,1,2-tetrachloroethane	13.564	131	82762	51.21	ug/L	97
98) ethylbenzene	13.559	91	302655	45.04	ug/L	100
99) m,p-xylene	13.669	106	235388	93.71	ug/L	97
100) o-xylene	14.083	106	123597	49.06	ug/L	92
101) styrene	14.094	104	201809	50.35	ug/L	96
103) bromoform	14.361	173	59726	51.94	ug/L	99
105) isopropylbenzene	14.424	105	327117	46.32	ug/L	99
107) bromobenzene	14.823	156	92715	48.15	ug/L	92
108) cyclohexanone	14.587	55	41632	221.03	ug/L	98
109) 1,1,2,2-tetrachloroethane	14.728	83	87641	43.19	ug/L	100
110) trans-1,4-dichloro-2-b...	14.770	53	19328	41.57	ug/L	97
111) 1,2,3-trichloropropane	14.802	110	19995	45.28	ug/L	93
112) n-propylbenzene	14.833	91	351383	44.48	ug/L	99
114) 2-chlorotoluene	14.980	126	79829	47.06	ug/L	92
115) 4-chlorotoluene	15.080	126	80891	48.06	ug/L	97
116) 1,3,5-trimethylbenzene	14.985	105	277483	46.56	ug/L	98
117) tert-butylbenzene	15.331	119	234010	46.62	ug/L	98
118) pentachloroethane	15.415	167	62013	48.47	ug/L	96
119) 1,2,4-trimethylbenzene	15.379	105	274605	45.91	ug/L	99
120) sec-butylbenzene	15.546	105	371795	47.35	ug/L	98
121) 1,3-dichlorobenzene	15.735	146	178588	46.51	ug/L	99
122) p-isopropyltoluene	15.662	119	317327	46.95	ug/L	98
123) 1,4-dichlorobenzene	15.814	146	182996	45.99	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4810-4810\
 Data File : 3d112242.d
 Acq On : 29 Jul 2015 11:25 pm
 Operator : ximenac
 Sample : ICC4810-50
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Jul 30 11:52:18 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
124) 1,2-dichlorobenzene	16.191	146	190124	46.14	ug/L	99
126) n-butylbenzene	16.065	92	162816	46.44	ug/L	100
128) 1,2-dibromo-3-chloropr...	16.925	157	26104	48.23	ug/L	96
129) 1,3,5-trichlorobenzene	17.093	180	206822	48.85	ug/L	98
130) 1,2,4-trichlorobenzene	17.659	180	213293	48.80	ug/L	99
131) hexachlorobutadiene	17.759	225	96893	47.19	ug/L	99
132) naphthalene	17.900	128	430037	44.69	ug/L	99
133) 1,2,3-trichlorobenzene	18.110	180	216874	47.93	ug/L	98
134) hexachloroethane	16.443	119	65619	56.94	ug/L	97
135) Benzyl chloride	15.924	91	165957	44.35	ug/L	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

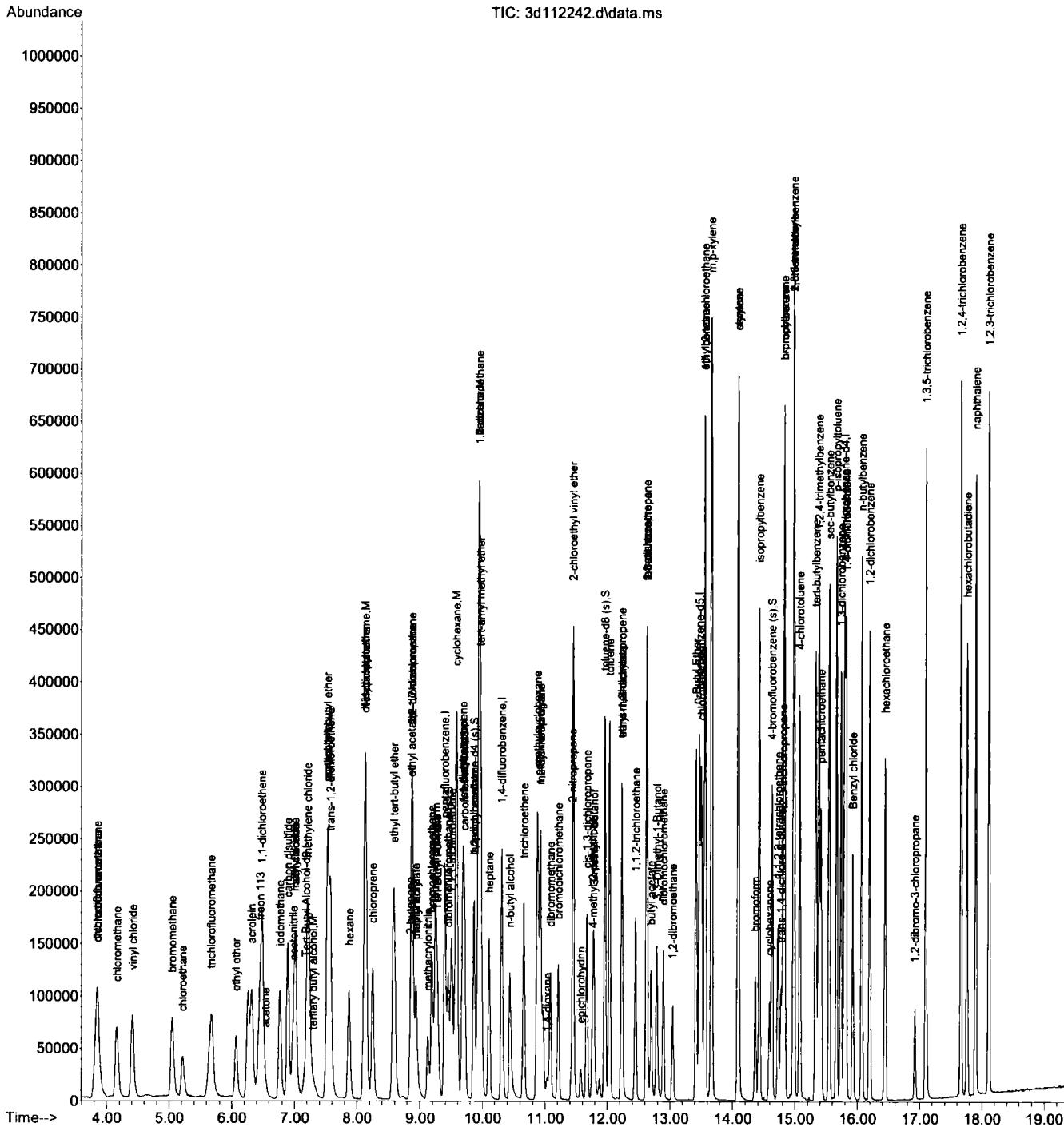
7.6.8



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112242.d
Acq On : 29 Jul 2015 11:25 pm
Operator : ximenac
Sample : ICC4810-50
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Jul 30 11:52:18 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration



7.6.3

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112243.d
 Acq On : 29 Jul 2015 11:53 pm
 Operator : ximenac
 Sample : IC4810-100
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 28 Sample Multiplier: 1

Quant Time: Jul 30 11:54:06 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.178	65	101373	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	208084	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	231990	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	189712	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	115891	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	80199	44.56	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	89.12%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	72486	39.01	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	78.02%	
80) toluene-d8 (s)	11.960	98	253534	51.32	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	102.64%	
106) 4-bromofluorobenzene (s)	14.629	95	90865	46.33	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	92.66%	
Target Compounds						
2) 1,4-dioxane	11.032	88	44544	2632.87	ug/L	88
3) tertiary butyl alcohol	7.288	59	108317	475.19	ug/L	98
8) chlorodifluoromethane	3.844	51	290111	70.40	ug/L	98
9) dichlorodifluoromethane	3.833	85	369475	87.57	ug/L	96
10) chloromethane	4.163	50	286392	92.30	ug/L	98
11) vinyl chloride	4.415	62	326335	96.28	ug/L	99
12) bromomethane	5.044	94	183337	109.84	ug/L	96
13) chloroethane	5.212	64	117683	93.03	ug/L	96
16) trichlorofluoromethane	5.668	101	322274	94.92	ug/L	99
18) ethyl ether	6.067	74	74650	89.85	ug/L	96
21) acrolein	6.329	56	6378	15.80	ug/L	93
22) 1,1-dichloroethene	6.481	61	249495	68.63	ug/L	94
23) acetone	6.533	43	45160	76.01	ug/L	96
24) allyl chloride	7.010	76	78943	75.00	ug/L #	86
25) acetonitrile	6.979	40	156090	683.52	ug/L	86
27) iodomethane	6.764	142	370642	85.95	ug/L	97
28) iso-butyl alcohol	9.684	41	53395	722.48	ug/L	96
29) carbon disulfide	6.890	76	633150	71.37	ug/L	98
30) methylene chloride	7.204	84	182921	74.55	ug/L	92
31) methyl acetate	6.995	43	106630	70.74	ug/L	95
32) methyl tert butyl ether	7.529	73	554719	81.54	ug/L	97
33) trans-1,2-dichloroethene	7.571	61	225371	70.65	ug/L	92
34) di-isopropyl ether	8.122	45	567957	73.71	ug/L	87
35) ethyl tert-butyl ether	8.583	59	584396	78.04	ug/L	98
36) 2-butanone	8.840	72	16056	93.61	ug/L	70
37) 1,1-dichloroethane	8.138	63	277583	74.24	ug/L	98
38) chloroprene	8.248	53	201905	71.97	ug/L	95
39) acrylonitrile	7.524	53	289356	422.60	ug/L	99
40) vinyl acetate	8.122	86	22420	93.45	ug/L	65
41) ethyl acetate	8.861	45	18603	81.89	ug/L	73
42) 2,2-dichloropropane	8.877	77	267502	76.99	ug/L	99
43) cis-1,2-dichloroethene	8.877	96	173618	80.42	ug/L	90
44) propionitrile	8.935	54	213487	841.82	ug/L	92
45) methyl acrylate	8.940	55	145849	87.46	ug/L	96
46) bromochloromethane	9.186	128	91593	90.37	ug/L #	86
47) tetrahydrofuran	9.223	42	48461	70.42	ug/L	95
48) chloroform	9.244	83	262330	76.63	ug/L	97
49) Tert-Butyl Formate	9.270	59	162008	83.88	ug/L	98
52) freon 113	6.455	151	150661	85.00	ug/L	89
53) methacrylonitrile	9.118	41	80998	78.79	ug/L	94
54) 1,1,1-trichloroethane	9.501	97	273762	84.95	ug/L	97

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112243.d
 Acq On : 29 Jul 2015 11:53 pm
 Operator : ximenac
 Sample : IC4810-100
 Misc : MS88759,V3D4810,5,,,.1
 ALS Vial : 28 Sample Multiplier: 1

Quant Time: Jul 30 11:54:06 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
57) tert-amyl methyl ether	9.973	73	548551	77.85	ug/L	97
59) epichlorohydrin	11.572	57	61959	474.88	ug/L	95
60) n-butyl alcohol	10.434	56	204395	5496.15	ug/L	92
61) cyclohexane	9.569	84	273210	96.33	ug/L	83
62) carbon tetrachloride	9.700	117	245891	102.20	ug/L	100
63) 1,1-dichloropropene	9.674	75	179730	84.51	ug/L	95
64) hexane	7.870	57	142993	72.43	ug/L	98
65) benzene	9.936	78	557400	85.59	ug/L	97
66) heptane	10.109	57	82958	79.32	ug/L	95
67) isopropyl acetate	9.857	43	258670	85.17	ug/L	98
68) 1,2-dichloroethane	9.952	62	168762	81.52	ug/L	98
69) trichloroethene	10.654	95	132679	91.70	ug/L	91
72) 2-nitropropane	11.430	41	48762	85.26	ug/L	96
73) 2-chloroethyl vinyl ether	11.446	63	440900	508.65	ug/L	99
74) methyl methacrylate	10.922	41	206219	87.44	ug/L	88
75) 1,2-dichloropropane	10.916	63	145208	86.27	ug/L	99
76) methylcyclohexane	10.880	83	269475	87.37	ug/L	94
77) dibromomethane	11.084	93	88547	93.29	ug/L	94
78) bromodichloromethane	11.210	83	179872	89.99	ug/L	97
79) cis-1,3-dichloropropene	11.671	75	220654	98.41	ug/L	92
81) 4-methyl-2-pentanone	11.761	58	56507	102.53	ug/L	97
82) toluene	12.033	92	322489	93.39	ug/L	97
83) 3-methyl-1-butanol	11.782	70	78706	2176.99	ug/L	88
84) trans-1,3-dichloropropene	12.232	75	199141	99.39	ug/L	91
85) ethyl methacrylate	12.222	69	172834	111.87	ug/L	97
86) 1,1,2-trichloroethane	12.447	83	103157	96.69	ug/L	98
87) 2-hexanone	12.620	58	47662	108.18	ug/L	97
89) tetrachloroethene	12.626	166	163840	105.46	ug/L	98
90) 1,3-dichloropropane	12.626	76	190824	94.99	ug/L	98
91) butyl acetate	12.694	56	84118	97.96	ug/L	92
92) 3,3-Dimethyl-1-Butanol	12.788	57	185997	1028.81	ug/L	99
93) dibromochloromethane	12.893	129	158524	103.96	ug/L	96
94) 1,2-dibromoethane	13.045	107	129781	97.22	ug/L	100
95) n-Butyl Ether	13.423	57	573195	90.93	ug/L	98
96) chlorobenzene	13.507	112	361557	93.35	ug/L	96
97) 1,1,2,2-tetrachloroethane	13.564	131	165727	104.49	ug/L	98
98) ethylbenzene	13.559	91	599961	90.98	ug/L	100
99) m,p-xylene	13.664	106	464345	188.35	ug/L	99
100) o-xylene	14.083	106	244273	98.79	ug/L	91
101) styrene	14.094	104	397073	100.93	ug/L	97
103) bromoform	14.361	173	124093	109.95	ug/L	98
105) isopropylbenzene	14.424	105	647770	96.66	ug/L	99
107) bromobenzene	14.823	156	180039	98.52	ug/L	91
108) cyclohexanone	14.587	55	221967	1241.78	ug/L	98
109) 1,1,2,2-tetrachloroethane	14.728	83	178365	92.62	ug/L	100
110) trans-1,4-dichloro-2-b...	14.770	53	40127	90.93	ug/L	90
111) 1,2,3-trichloropropane	14.802	110	40860	97.49	ug/L	97
112) n-propylbenzene	14.833	91	689129	91.93	ug/L	100
114) 2-chlorotoluene	14.980	126	159391	99.02	ug/L	93
115) 4-chlorotoluene	15.079	126	155133	97.11	ug/L	96
116) 1,3,5-trimethylbenzene	14.985	105	556138	98.32	ug/L	97
117) tert-butylbenzene	15.331	119	477104	100.17	ug/L	98
118) pentachloroethane	15.415	167	133536	109.97	ug/L	96
119) 1,2,4-trimethylbenzene	15.378	105	547107	96.38	ug/L	98
120) sec-butylbenzene	15.546	105	751437	100.84	ug/L	98
121) 1,3-dichlorobenzene	15.735	146	344497	94.55	ug/L	99
122) p-isopropyltoluene	15.661	119	639663	99.74	ug/L	98
123) 1,4-dichlorobenzene	15.813	146	351034	92.96	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112243.d
 Acq On : 29 Jul 2015 11:53 pm
 Operator : ximenac
 Sample : IC4810-100
 Misc : MS88759,V3D4810,5,,,,1
 ALS Vial : 28 Sample Multiplier: 1

Quant Time: Jul 30 11:54:06 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.191	146	368245	94.17	ug/L	98
126) n-butylbenzene	16.065	92	316811	95.22	ug/L	98
128) 1,2-dibromo-3-chloropr...	16.920	157	55537	108.12	ug/L	97
129) 1,3,5-trichlorobenzene	17.093	180	405076	100.83	ug/L	99
130) 1,2,4-trichlorobenzene	17.659	180	423668	102.13	ug/L	99
131) hexachlorobutadiene	17.759	225	199123	102.18	ug/L	99
132) naphthalene	17.895	128	893950	97.89	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	434998	101.30	ug/L	98
134) hexachloroethane	16.443	119	142757	130.55	ug/L	97
135) Benzyl chloride	15.924	91	325837	91.75	ug/L	97

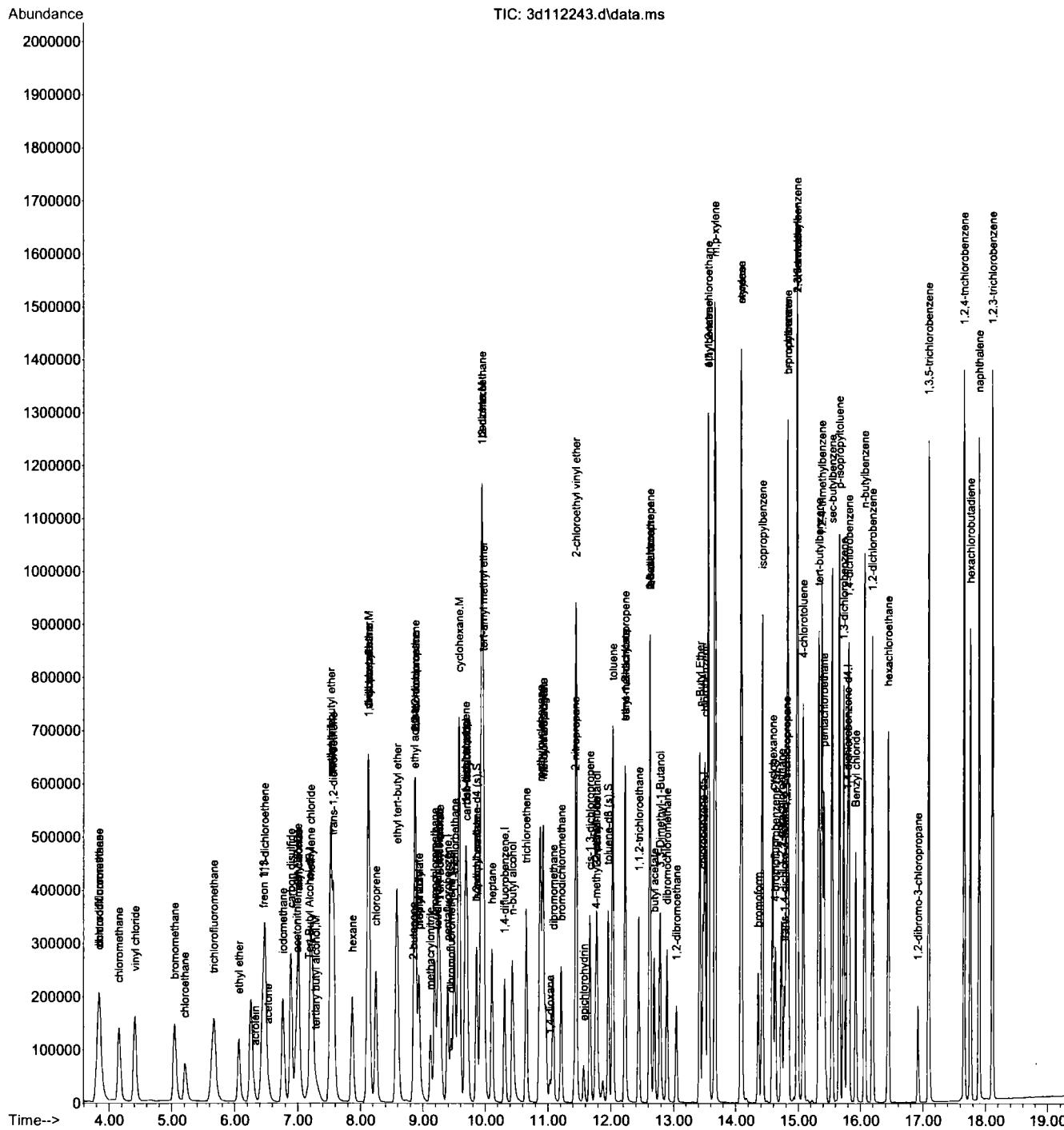
(#) = qualifier out of range (m) = manual integration (+) = signals summed

7.6.9

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112243.d
Acq On : 29 Jul 2015 11:53 pm
Operator : ximenac
Sample : IC4810-100
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 28 Sample Multiplier: 1

Quant Time: Jul 30 11:54:06 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Jul 30 08:37:34 2015
Response via : Initial Calibration



M3D4810.M Thu Aug 06 13:50:25 2015 3D

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LABORATORIES
JC1106

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112244.d
 Acq On : 30 Jul 2015 12:20 am
 Operator : ximenac
 Sample : IC4810-200
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 29 Sample Multiplier: 1

Quant Time: Jul 30 11:55:56 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.183	65	113652	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	211907	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	243867	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	189787	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	123773	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	84139	45.91	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	91.82%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	75693	40.00	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	80.00%	
80) toluene-d8 (s)	11.965	98	260601	50.18	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	100.36%	
106) 4-bromofluorobenzene (s)	14.629	95	90862	43.38	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	86.76%	
Target Compounds						
				Qvalue		
2) 1,4-dioxane	11.032	88	95700	5045.42	ug/L	84
3) tertiary butyl alcohol	7.288	59	235989	923.44	ug/L	96
8) chlorodifluoromethane	3.844	51	590213	140.64	ug/L	97
9) dichlorodifluoromethane	3.838	85	728302	169.51	ug/L	99
10) chloromethane	4.169	50	622229	196.91	ug/L	98
11) vinyl chloride	4.420	62	703897	203.94	ug/L	99
12) bromomethane	5.044	94	357394	210.27	ug/L	99
13) chloroethane	5.207	64	234170	181.77	ug/L	97
16) trichlorofluoromethane	5.673	101	651329	188.38	ug/L	100
18) ethyl ether	6.067	74	153481	181.39	ug/L	99
22) 1,1-dichloroethene	6.481	61	512563	138.46	ug/L	94
23) acetone	6.539	43	92034	152.11	ug/L	95
24) allyl chloride	7.010	76	161550	150.72	ug/L	# 84
25) acetonitrile	6.969	40	323563	1391.33	ug/L	97
27) iodomethane	6.764	142	767087	174.67	ug/L	98
28) iso-butyl alcohol	9.684	41	118343	1572.40	ug/L	91
29) carbon disulfide	6.890	76	1299931	143.88	ug/L	98
30) methylene chloride	7.204	84	373651	149.54	ug/L	93
31) methyl acetate	6.995	43	228717	148.99	ug/L	96
32) methyl tert butyl ether	7.530	73	1138567	164.33	ug/L	97
33) trans-1,2-dichloroethene	7.571	61	461296	142.00	ug/L	92
34) di-isopropyl ether	8.122	45	1165663	148.55	ug/L	84
35) ethyl tert-butyl ether	8.589	59	1209427	158.60	ug/L	97
36) 2-butanone	8.840	72	34182	195.70	ug/L	70
37) 1,1-dichloroethane	8.138	63	566169	148.70	ug/L	99
38) chloroprene	8.248	53	413880	144.88	ug/L	94
39) acrylonitrile	7.524	53	602572	864.16	ug/L	98
40) vinyl acetate	8.122	86	47648	195.02	ug/L	74
41) ethyl acetate	8.856	45	38508	166.45	ug/L	69
42) 2,2-dichloropropane	8.877	77	537523	151.91	ug/L	98
43) cis-1,2-dichloroethene	8.877	96	358567	163.08	ug/L	90
44) propionitrile	8.935	54	460126	1781.63	ug/L	91
45) methyl acrylate	8.935	55	311809	183.62	ug/L	97
46) bromochloromethane	9.186	128	188533	182.66	ug/L	# 86
47) tetrahydrofuran	9.228	42	105272	150.22	ug/L	94
48) chloroform	9.244	83	542991	155.76	ug/L	98
49) Tert-Butyl Formate	9.275	59	343297	174.54	ug/L	98
52) freon 113	6.449	151	296095	164.03	ug/L	88
53) methacrylonitrile	9.118	41	174920	167.08	ug/L	93
54) 1,1,1-trichloroethane	9.496	97	561095	170.96	ug/L	97
57) tert-amyl methyl ether	9.973	73	1153060	160.70	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112244.d
 Acq On : 30 Jul 2015 12:20 am
 Operator : ximenac
 Sample : IC4810-200
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 29 Sample Multiplier: 1

Quant Time: Jul 30 11:55:56 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
59) epichlorohydrin	11.572	57	130867	954.18	ug/L	94
60) n-butyl alcohol	10.434	56	459085	11743.50	ug/L	91
61) cyclohexane	9.569	84	548425	183.94	ug/L	85
62) carbon tetrachloride	9.705	117	498383	197.06	ug/L	99
63) 1,1-dichloropropene	9.674	75	375601	168.01	ug/L	94
64) hexane	7.870	57	286276	137.95	ug/L	95
65) benzene	9.936	78	1180256	172.41	ug/L	98
66) heptane	10.109	57	166976	151.87	ug/L	96
67) isopropyl acetate	9.857	43	556260	174.23	ug/L	99
68) 1,2-dichloroethane	9.952	62	355113	163.19	ug/L	99
69) trichloroethene	10.654	95	282796	185.94	ug/L	93
72) 2-nitropropane	11.425	41	99101	164.84	ug/L	98
73) 2-chloroethyl vinyl ether	11.451	63	883990	970.16	ug/L	97
74) methyl methacrylate	10.922	41	434514	175.27	ug/L	87
75) 1,2-dichloropropane	10.917	63	303367	171.45	ug/L	98
76) methylcyclohexane	10.875	83	554075	170.89	ug/L	93
77) dibromomethane	11.079	93	185760	186.18	ug/L	94
78) bromodichloromethane	11.210	83	385984	183.71	ug/L	96
79) cis-1,3-dichloropropene	11.672	75	451689	191.64	ug/L	93
81) 4-methyl-2-pentanone	11.761	58	123218	212.69	ug/L	94
82) toluene	12.033	92	663934	182.90	ug/L	98
83) 3-methyl-1-butanol	11.782	70	180645	4753.24	ug/L	87
84) trans-1,3-dichloropropene	12.233	75	404138	191.88	ug/L	91
85) ethyl methacrylate	12.222	69	354848	218.49	ug/L	97
86) 1,1,2-trichloroethane	12.448	83	211819	188.88	ug/L	98
87) 2-hexanone	12.615	58	101312	218.76	ug/L	97
89) tetrachloroethene	12.626	166	330772	212.82	ug/L	98
90) 1,3-dichloropropane	12.626	76	381982	190.08	ug/L	98
91) butyl acetate	12.694	56	177748	206.92	ug/L	95
92) 3,3-Dimethyl-1-Butanol	12.788	57	447144	2472.31	ug/L	99
93) dibromochloromethane	12.893	129	323679	212.19	ug/L	96
94) 1,2-dibromoethane	13.045	107	263194	197.08	ug/L	99
95) n-Butyl Ether	13.423	57	1137615	180.39	ug/L	97
96) chlorobenzene	13.507	112	729983	188.41	ug/L	96
97) 1,1,1,2-tetrachloroethane	13.564	131	356803	224.87	ug/L	97
98) ethylbenzene	13.559	91	1207257	182.99	ug/L	99
99) m,p-xylene	13.669	106	932385	378.06	ug/L	99
100) o-xylene	14.083	106	501975	202.94	ug/L	94
101) styrene	14.094	104	798986	203.01	ug/L	98
103) bromoform	14.356	173	257021	227.64	ug/L	98
105) isopropylbenzene	14.424	105	1349790	188.60	ug/L	99
107) bromobenzene	14.823	156	362575	185.78	ug/L	91
108) cyclohexanone	14.587	55	459103	2404.85	ug/L	97
109) 1,1,2,2-tetrachloroethane	14.728	83	373746	181.72	ug/L	100
110) trans-1,4-dichloro-2-b...	14.770	53	84850	180.04	ug/L	89
111) 1,2,3-trichloropropene	14.802	110	84031	187.73	ug/L	99
112) n-propylbenzene	14.833	91	1372797	171.47	ug/L	99
114) 2-chlorotoluene	14.980	126	321057	186.74	ug/L	95
115) 4-chlorotoluene	15.079	126	315361	184.85	ug/L	95
116) 1,3,5-trimethylbenzene	14.985	105	1167012	193.18	ug/L	97
117) tert-butylbenzene	15.331	119	1064850	209.33	ug/L	98
118) pentachloroethane	15.415	167	305312	235.43	ug/L	94
119) 1,2,4-trimethylbenzene	15.378	105	1127745	186.01	ug/L	98
120) sec-butylbenzene	15.546	105	1595696	200.50	ug/L	99
121) 1,3-dichlorobenzene	15.735	146	728618	187.24	ug/L	98
122) p-isopropyltoluene	15.661	119	1336685	195.15	ug/L	98
123) 1,4-dichlorobenzene	15.814	146	752826	186.66	ug/L	99
124) 1,2-dichlorobenzene	16.191	146	797169	190.87	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\V3D4810-4810\
 Data File : 3d112244.d
 Acq On : 30 Jul 2015 12:20 am
 Operator : ximenac
 Sample : IC4810-200
 Misc : MS88759, V3D4810,5,,,1
 ALS Vial : 29 Sample Multiplier: 1

Quant Time: Jul 30 11:55:56 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
126) n-butylbenzene	16.065	92	663120	186.61	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.920	157	128163	233.62	ug/L	97
129) 1,3,5-trichlorobenzene	17.093	180	925275	215.65	ug/L	98
130) 1,2,4-trichlorobenzene	17.659	180	958374	216.32	ug/L	99
131) hexachlorobutadiene	17.759	225	461755	221.87	ug/L	99
132) naphthalene	17.900	128	1999333	204.98	ug/L	99
133) 1,2,3-trichlorobenzene	18.110	180	948626	206.85	ug/L	99
134) hexachloroethane	16.443	119	334395	286.32	ug/L	97
135) Benzyl chloride	15.924	91	717668	189.21	ug/L	97

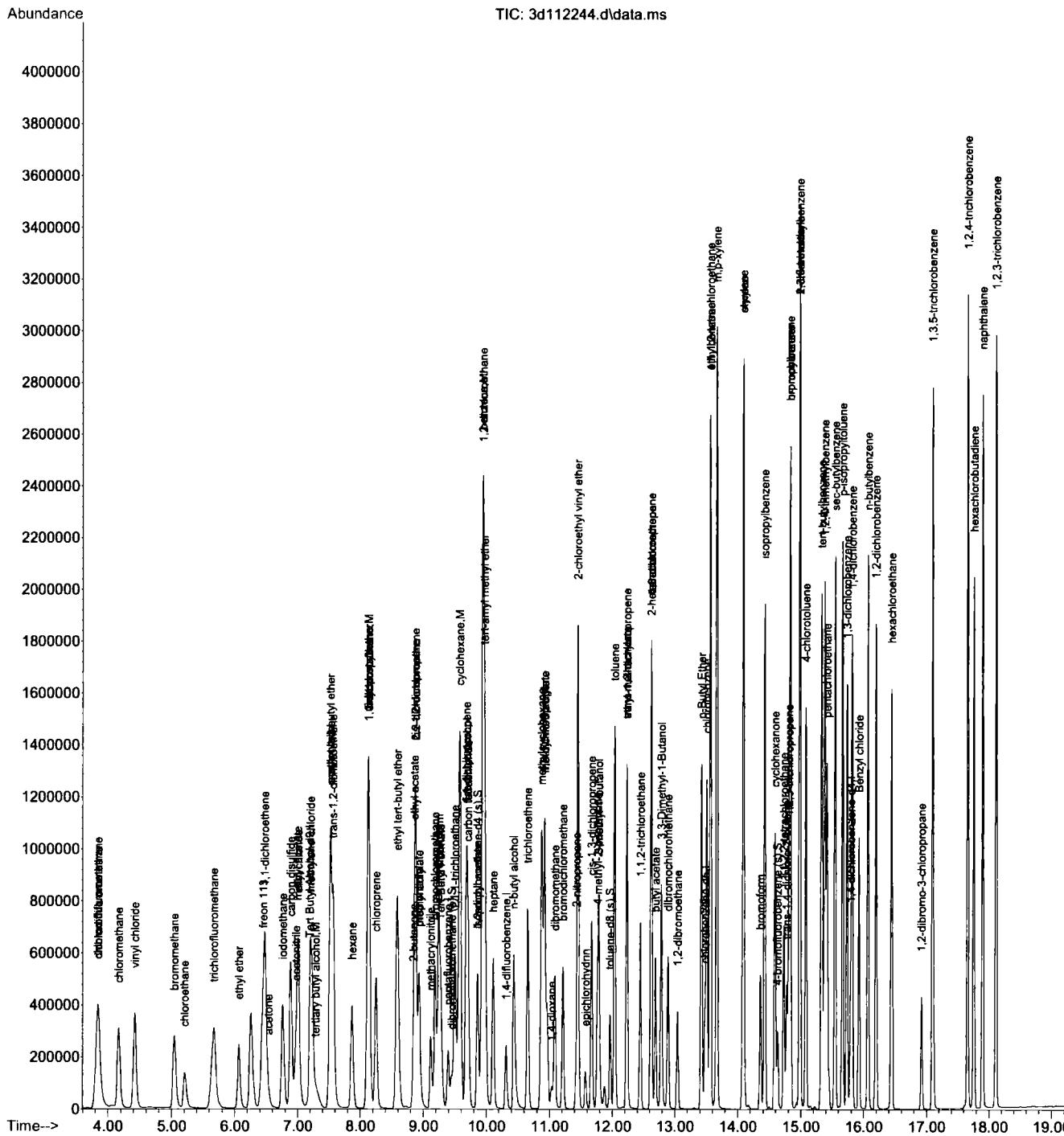
(#) = qualifier out of range (m) = manual integration (+) = signals summed

7.6.10
7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112244.d
 Acq On : 30 Jul 2015 12:20 am
 Operator : ximenac
 Sample : IC4810-200
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 29 Sample Multiplier: 1

Quant Time: Jul 30 11:55:56 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Jul 30 08:37:34 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112247.d
 Acq On : 30 Jul 2015 1:41 am
 Operator : ximenac
 Sample : ICV4810-50
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 32 Sample Multiplier: 1

Quant Time: Aug 06 13:51:00 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) Tert Butyl Alcohol-d9	7.178	65	98586	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	207113	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	228341	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	187656	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	123839	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	80387	50.66	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	101.32%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	71580	48.69	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	97.38%	
80) toluene-d8 (s)	11.960	98	251682	50.33	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	100.66%	
106) 4-bromofluorobenzene (s)	14.629	95	94440	49.57	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	99.14%	
Target Compounds						
2) 1,4-dioxane	11.037	88	22005	1381.38	ug/L	98
3) tertiary butyl alcohol	7.288	59	53644	251.95	ug/L	96
8) chlorodifluoromethane	3.854	51	117664	41.85	ug/L	97
9) dichlorodifluoromethane	3.849	85	139646	37.10	ug/L	98
10) chloromethane	4.164	50	124313	41.53	ug/L	99
11) vinyl chloride	4.410	62	136816	45.57	ug/L	100
12) bromomethane	5.050	94	89616	46.20	ug/L	97
13) chloroethane	5.212	64	65378	51.42	ug/L	99
16) trichlorofluoromethane	5.679	101	135791	43.34	ug/L	99
18) ethyl ether	6.067	74	36422	47.07	ug/L	95
21) acrolein	6.318	56	157273	506.30	ug/L	99
22) 1,1-dichloroethene	6.491	61	103370	41.96	ug/L	96
23) acetone	6.539	43	22416	48.30	ug/L	87
24) allyl chloride	7.016	76	41133	53.76	ug/L	98
25) acetonitrile	6.974	40	78456	508.42	ug/L	93
27) iodomethane	6.764	142	170305	48.62	ug/L	98
28) iso-butyl alcohol	9.684	41	25600	496.49	ug/L	91
29) carbon disulfide	6.890	76	280190	44.98	ug/L	99
30) methylene chloride	7.205	84	88751	48.08	ug/L	99
31) methyl acetate	7.000	43	53457	51.85	ug/L	98
32) methyl tert butyl ether	7.530	73	536282	98.07	ug/L	94
33) trans-1,2-dichloroethene	7.577	61	101915	45.34	ug/L	97
34) di-isopropyl ether	8.122	45	283901	49.11	ug/L	94
35) ethyl tert-butyl ether	8.589	59	290928	53.42	ug/L	97
36) 2-butanone	8.840	72	7736	52.38	ug/L	97
37) 1,1-dichloroethane	8.143	63	132194	49.50	ug/L	98
38) chloroprene	8.248	53	97195	49.73	ug/L	98
39) acrylonitrile	7.524	53	141741	265.20	ug/L	94
40) vinyl acetate	8.127	86	11484	57.88	ug/L	87
41) ethyl acetate	8.861	45	9823	48.35	ug/L	74
42) 2,2-dichloropropane	8.882	77	115296	41.26	ug/L	98
43) cis-1,2-dichloroethene	8.882	96	81068	44.91	ug/L	97
44) propionitrile	8.940	54	98024	492.17	ug/L	98
45) methyl acrylate	8.945	55	69484	51.75	ug/L	99
46) bromochloromethane	9.192	128	44037	50.70	ug/L	98
47) tetrahydrofuran	9.223	42	24949	47.69	ug/L	97
48) chloroform	9.244	83	125087	48.58	ug/L	100
49) Tert-Butyl Formate	9.275	59	81940	54.66	ug/L	# 99
52) freon 113	6.449	151	71327	47.03	ug/L	99
53) methacrylonitrile	9.123	41	39614	50.91	ug/L	98
54) 1,1,1-trichloroethane	9.501	97	123034	46.28	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
 Data File : 3d112247.d
 Acq On : 30 Jul 2015 1:41 am
 Operator : ximenac
 Sample : ICV4810-50
 Misc : MS88759, V3D4810, 5, , , 1
 ALS Vial : 32 Sample Multiplier: 1

Quant Time: Aug 06 13:51:00 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
57) tert-amyl methyl ether	9.973	73	272329	50.03	ug/L	99
59) epichlorohydrin	11.572	57	31673	252.67	ug/L	98
60) n-butyl alcohol	10.439	56	99738	2675.22	ug/L	98
61) cyclohexane	9.574	84	116852	43.70	ug/L	99
62) carbon tetrachloride	9.705	117	107740	48.79	ug/L	96
63) 1,1-dichloropropene	9.679	75	80084	48.69	ug/L	98
64) hexane	7.870	57	47903	32.61	ug/L	97
65) benzene	9.936	78	256456	47.80	ug/L	99
66) heptane	10.104	57	37301	44.70	ug/L	95
67) isopropyl acetate	9.857	43	142485	61.05	ug/L	96
68) 1,2-dichloroethane	9.952	62	82754	51.64	ug/L	97
69) trichloroethene	10.660	95	61075	47.08	ug/L	97
72) 2-nitropropane	11.430	41	24588	51.85	ug/L	100
73) 2-chloroethyl vinyl ether	11.451	63	248238	301.43	ug/L	98
74) methyl methacrylate	10.922	41	100489	54.78	ug/L	96
75) 1,2-dichloropropane	10.922	63	70280	50.11	ug/L	99
76) methylcyclohexane	10.875	83	124476	51.30	ug/L	97
77) dibromomethane	11.084	93	42338	51.31	ug/L	95
78) bromodichloromethane	11.210	83	86540	50.71	ug/L	99
79) cis-1,3-dichloropropene	11.672	75	105521	52.67	ug/L	97
81) 4-methyl-2-pentanone	11.761	58	27358	53.59	ug/L	98
82) toluene	12.033	92	145656	49.40	ug/L	100
83) 3-methyl-1-butanol	11.787	70	37805	979.95	ug/L	96
84) trans-1,3-dichloropropene	12.233	75	90548	50.86	ug/L	98
85) ethyl methacrylate	12.227	69	79460	50.50	ug/L	97
86) 1,1,2-trichloroethane	12.448	83	49718	50.20	ug/L	97
87) 2-hexanone	12.621	58	21853	52.36	ug/L	97
89) tetrachloroethene	12.626	166	83541	52.29	ug/L	99
90) 1,3-dichloropropane	12.631	76	90638	49.78	ug/L	99
91) butyl acetate	12.694	56	41482	55.05	ug/L	98
92) 3,3-Dimethyl-1-Butanol	12.788	57	77128	428.24	ug/L	99
93) dibromochloromethane	12.893	129	74597	52.54	ug/L	98
94) 1,2-dibromoethane	13.045	107	62203	51.95	ug/L	99
95) n-Butyl Ether	13.423	57	265121	46.70	ug/L	99
96) chlorobenzene	13.507	112	171846	51.06	ug/L	99
97) 1,1,1,2-tetrachloroethane	13.564	131	75616	50.67	ug/L	98
98) ethylbenzene	13.559	91	273475	47.44	ug/L	99
99) m,p-xylene	13.669	106	214664	99.59	ug/L	97
100) o-xylene	14.083	106	114530	52.02	ug/L	99
101) styrene	14.094	104	192022	51.83	ug/L	100
103) bromoform	14.356	173	60528	54.64	ug/L	99
105) isopropylbenzene	14.424	105	289437	46.57	ug/L	99
107) bromobenzene	14.823	156	89866	48.99	ug/L	98
108) cyclohexanone	14.587	55	101056	487.84	ug/L	97
109) 1,1,2,2-tetrachloroethane	14.728	83	87322	48.55	ug/L	99
110) trans-1,4-dichloro-2-b...	14.770	53	19558	49.14	ug/L	97
111) 1,2,3-trichloropropane	14.802	110	20269	49.35	ug/L	98
112) n-propylbenzene	14.833	91	336151	49.68	ug/L	99
114) 2-chlorotoluene	14.980	126	75205	47.58	ug/L	99
115) 4-chlorotoluene	15.080	126	77133	50.26	ug/L	93
116) 1,3,5-trimethylbenzene	14.985	105	251941	47.64	ug/L	97
117) tert-butylbenzene	15.331	119	208377	47.82	ug/L	98
118) pentachloroethane	15.415	167	49840	41.94	ug/L	97
119) 1,2,4-trimethylbenzene	15.378	105	268403	50.55	ug/L	99
120) sec-butylbenzene	15.546	105	329100	47.26	ug/L	99
121) 1,3-dichlorobenzene	15.735	146	174542	51.18	ug/L	98
122) p-isopropyltoluene	15.662	119	293093	48.20	ug/L	98
123) 1,4-dichlorobenzene	15.814	146	181410	49.58	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4810-4810\
 Data File : 3d112247.d
 Acq On : 30 Jul 2015 1:41 am
 Operator : ximenac
 Sample : ICV4810-50
 Misc : MS88759,V3D4810,5,,,1
 ALS Vial : 32 Sample Multiplier: 1

Quant Time: Aug 06 13:51:00 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

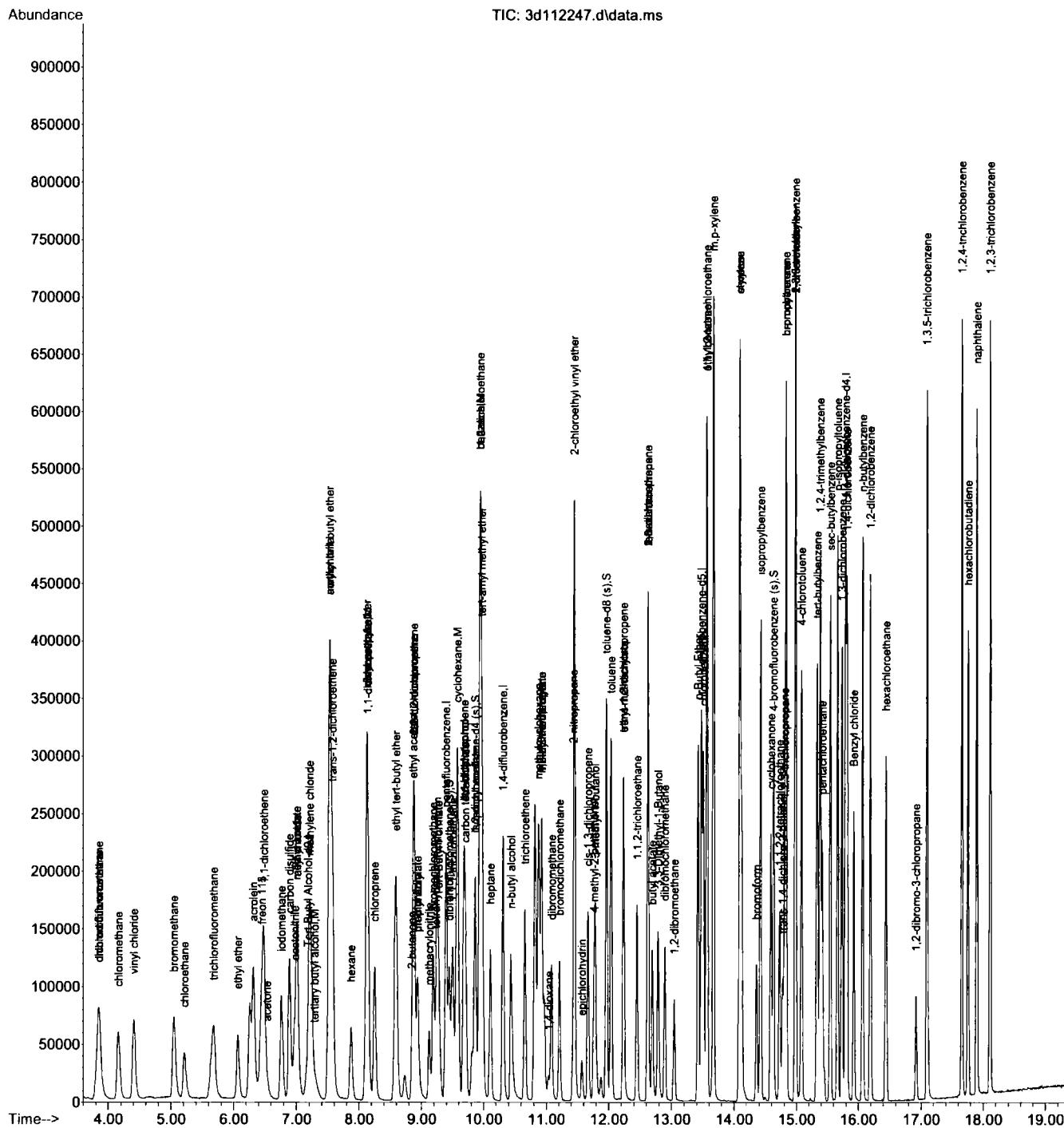
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.191	146	192407	52.33	ug/L	100
126) n-butylbenzene	16.065	92	153250	48.17	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.925	157	27972	52.06	ug/L	94
129) 1,3,5-trichlorobenzene	17.093	180	203232	52.06	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	210402	52.43	ug/L	98
131) hexachlorobutadiene	17.759	225	88768	44.69	ug/L	99
132) naphthalene	17.900	128	440270	51.80	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	213504	51.25	ug/L	99
134) hexachloroethane	16.443	119	59302	47.22	ug/L	99
135) Benzyl chloride	15.924	91	175556	49.14	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\v3d4801-4810\
Data File : 3d112247.d
Acq On : 30 Jul 2015 1:41 am
Operator : ximenac
Sample : ICV4810-50
Misc : MS88759,V3D4810,5,,,1
ALS Vial : 32 Sample Multiplier: 1

Quant Time: Aug 06 13:51:00 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112517.D
 Acq On : 11 Aug 2015 11:16 am
 Operator : ximenac
 Sample : cc4810-20
 Misc : MS89341,V3D4822,5,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 11 16:10:05 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.162	65	105534	500.00	ug/L	-0.02
4) pentafluorobenzene	9.391	168	204849	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.314	114	234799	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	194668	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.793	152	122430	50.00	ug/L	0.00

System Monitoring Compounds

50) dibromofluoromethane (s)	9.443	113	84380	53.76	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 107.52%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	77968	53.63	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 107.26%	
80) toluene-d8 (s)	11.960	98	258676	50.30	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 100.60%	
106) 4-bromofluorobenzene (s)	14.634	95	94609	50.23	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.46%	

Target Compounds

					Qvalue
2) 1,4-dioxane	11.037	88	9830	576.46	ug/L 88
3) tertiary butyl alcohol	7.299	59	22677	99.50	ug/L 94
8) chlorodifluoromethane	3.859	51	54443	19.58	ug/L 97
9) dichlorodifluoromethane	3.859	85	60799	16.33	ug/L 99
10) chloromethane	4.163	50	45185	15.26	ug/L 98
11) vinyl chloride	4.410	62	51634	17.39	ug/L 97
12) bromomethane	5.055	94	33348	17.38	ug/L 94
13) chloroethane	5.223	64	22579	17.95	ug/L 98
16) trichlorofluoromethane	5.684	101	63301	20.43	ug/L 99
18) ethyl ether	6.072	74	16113	21.05	ug/L 97
21) acrolein	6.324	56	58219	189.49	ug/L 95
22) 1,1-dichloroethene	6.486	61	49353	20.26	ug/L 96
23) acetone	6.539	43	11209	24.42	ug/L 85
24) allyl chloride	7.016	76	16135	21.32	ug/L 98
25) acetonitrile	6.979	40	35265	231.06	ug/L 90
27) iodomethane	6.764	142	77335	22.32	ug/L 95
28) iso-butyl alcohol	9.690	41	10690	209.61	ug/L 93
29) carbon disulfide	6.895	76	131042	21.27	ug/L 99
30) methylene chloride	7.199	84	40299	22.07	ug/L 97
31) methyl acetate	6.995	43	22642	22.20	ug/L 98
32) methyl tert butyl ether	7.535	73	113652	21.01	ug/L 97
33) trans-1,2-dichloroethene	7.577	61	47282	21.27	ug/L 99
34) di-isopropyl ether	8.122	45	102463	17.92	ug/L 95
35) ethyl tert-butyl ether	8.589	59	104578	19.41	ug/L 98
36) 2-butanone	8.851	72	2997	20.52	ug/L 86
37) 1,1-dichloroethane	8.143	63	57359	21.72	ug/L 95
38) chloroprene	8.248	53	33099	17.12	ug/L 96
39) acrylonitrile	7.529	53	62044	117.37	ug/L 96
40) vinyl acetate	8.132	86	4070	20.74	ug/L 58
41) ethyl acetate	8.866	45	3592	17.88	ug/L # 24
42) 2,2-dichloropropane	8.887	77	58652	21.22	ug/L 98
43) cis-1,2-dichloroethene	8.877	96	34480	19.31	ug/L 94
44) propionitrile	8.940	54	45348	230.20	ug/L 85
45) methyl acrylate	8.945	55	29328	22.09	ug/L 97
46) bromochloromethane	9.186	128	18951	22.06	ug/L 95
47) tetrahydrofuran	9.223	42	9958	19.24	ug/L 98
48) chloroform	9.244	83	53107	20.85	ug/L 100
49) Tert-Butyl Formate	9.281	59	28385	19.14	ug/L # 96
52) freon 113	6.460	151	30401	20.27	ug/L 93
53) methacrylonitrile	9.129	41	15091	19.61	ug/L 97
54) 1,1,1-trichloroethane	9.501	97	58473	22.24	ug/L 97

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112517.D
 Acq On : 11 Aug 2015 11:16 am
 Operator : ximenac
 Sample : cc4810-20
 Misc : MS89341,V3D4822,5,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 11 16:10:05 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
57) tert-amyl methyl ether	9.973	73	101825	18.91	ug/L	95
59) epichlorohydrin	11.577	57	12217	94.78	ug/L	96
60) n-butyl alcohol	10.445	56	38470	1003.48	ug/L	96
61) cyclohexane	9.574	84	54644	19.87	ug/L	95
62) carbon tetrachloride	9.705	117	52750	23.23	ug/L	96
63) 1,1-dichloropropene	9.674	75	34793	20.57	ug/L	97
64) hexane	7.870	57	25436	16.84	ug/L	99
65) benzene	9.936	78	112369	20.37	ug/L	99
66) heptane	10.109	57	16118	18.78	ug/L	97
67) isopropyl acetate	9.863	43	46424	19.34	ug/L	97
68) 1,2-dichloroethane	9.957	62	36200	21.97	ug/L	99
69) trichloroethene	10.660	95	26892	20.16	ug/L	99
72) 2-nitropropane	11.441	41	9394	19.26	ug/L	91
73) 2-chloroethyl vinyl ether	11.451	63	84918	100.28	ug/L	99
74) methyl methacrylate	10.927	41	38364	20.34	ug/L	91
75) 1,2-dichloropropane	10.922	63	29639	20.55	ug/L	97
76) methylcyclohexane	10.880	83	48923	19.61	ug/L	95
77) dibromomethane	11.079	93	19442	22.92	ug/L	91
78) bromodichloromethane	11.210	83	37345	21.28	ug/L	97
79) cis-1,3-dichloropropene	11.672	75	43256	21.00	ug/L	98
81) 4-methyl-2-pentanone	11.766	58	11169	21.27	ug/L	96
82) toluene	12.039	92	63714	21.01	ug/L	96
83) 3-methyl-1-butanol	11.787	70	15186	382.81	ug/L	98
84) trans-1,3-dichloropropene	12.238	75	39200	21.41	ug/L	95
85) ethyl methacrylate	12.227	69	29508	18.24	ug/L	96
86) 1,1,2-trichloroethane	12.447	83	21807	21.41	ug/L	96
87) 2-hexanone	12.626	58	8021	18.69	ug/L	83
89) tetrachloroethene	12.626	166	33370	20.13	ug/L	97
90) 1,3-dichloropropane	12.631	76	39846	21.10	ug/L	98
91) butyl acetate	12.699	56	14578	18.65	ug/L	92
92) 3,3-Dimethyl-1-Butanol	12.788	57	28725	153.74	ug/L	95
93) dibromochloromethane	12.898	129	32852	22.30	ug/L	98
94) 1,2-dibromoethane	13.045	107	26875	21.64	ug/L	96
95) n-Butyl Ether	13.423	57	99516	16.90	ug/L	97
96) chlorobenzene	13.507	112	73819	21.14	ug/L	98
97) 1,1,2,2-tetrachloroethane	13.564	131	33807	21.84	ug/L	99
98) ethylbenzene	13.564	91	117533	19.66	ug/L	97
99) m,p-xylene	13.669	106	91131	40.75	ug/L	99
100) o-xylene	14.089	106	47187	20.66	ug/L	94
101) styrene	14.099	104	72615	18.90	ug/L	98
103) bromoform	14.361	173	25629	22.30	ug/L	97
105) isopropylbenzene	14.429	105	121251	19.74	ug/L	98
107) bromobenzene	14.828	156	36987	20.40	ug/L	96
108) cyclohexanone	14.592	55	47206	259.91	ug/L	97
109) 1,1,2,2-tetrachloroethane	14.733	83	39443	22.18	ug/L	92
110) trans-1,4-dichloro-2-b...	14.775	53	7247	18.42	ug/L	96
111) 1,2,3-trichloropropane	14.807	110	9530	23.47	ug/L	90
112) n-propylbenzene	14.838	91	136729	20.44	ug/L	100
114) 2-chlorotoluene	14.985	126	32099	20.54	ug/L	97
115) 4-chlorotoluene	15.085	126	31828	20.98	ug/L	95
116) 1,3,5-trimethylbenzene	14.985	105	108453	20.74	ug/L	97
117) tert-butylbenzene	15.331	119	82994	19.26	ug/L	98
118) pentachloroethane	15.415	167	25378	21.60	ug/L	96
119) 1,2,4-trimethylbenzene	15.384	105	106276	20.25	ug/L	98
120) sec-butylbenzene	15.546	105	140165	20.36	ug/L	98
121) 1,3-dichlorobenzene	15.740	146	71994	21.35	ug/L	99
122) p-isopropyltoluene	15.667	119	120898	20.11	ug/L	99
123) 1,4-dichlorobenzene	15.819	146	75263	20.81	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112517.D
 Acq On : 11 Aug 2015 11:16 am
 Operator : ximenac
 Sample : cc4810-20
 Misc : MS89341,V3D4822,5,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 11 16:10:05 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.196	146	78353	21.56	ug/L	99
126) n-butylbenzene	16.070	92	63468	20.18	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.930	157	11422	21.50	ug/L	95
129) 1,3,5-trichlorobenzene	17.098	180	83689	21.69	ug/L	100
130) 1,2,4-trichlorobenzene	17.664	180	87613	22.09	ug/L	99
131) hexachlorobutadiene	17.764	225	39454	20.09	ug/L	97
132) naphthalene	17.900	128	185567	22.08	ug/L	99
133) 1,2,3-trichlorobenzene	18.115	180	92686	22.50	ug/L	99
134) hexachloroethane	16.443	119	24651	19.85	ug/L	99
135) Benzyl chloride	15.929	91	70680	20.01	ug/L	99

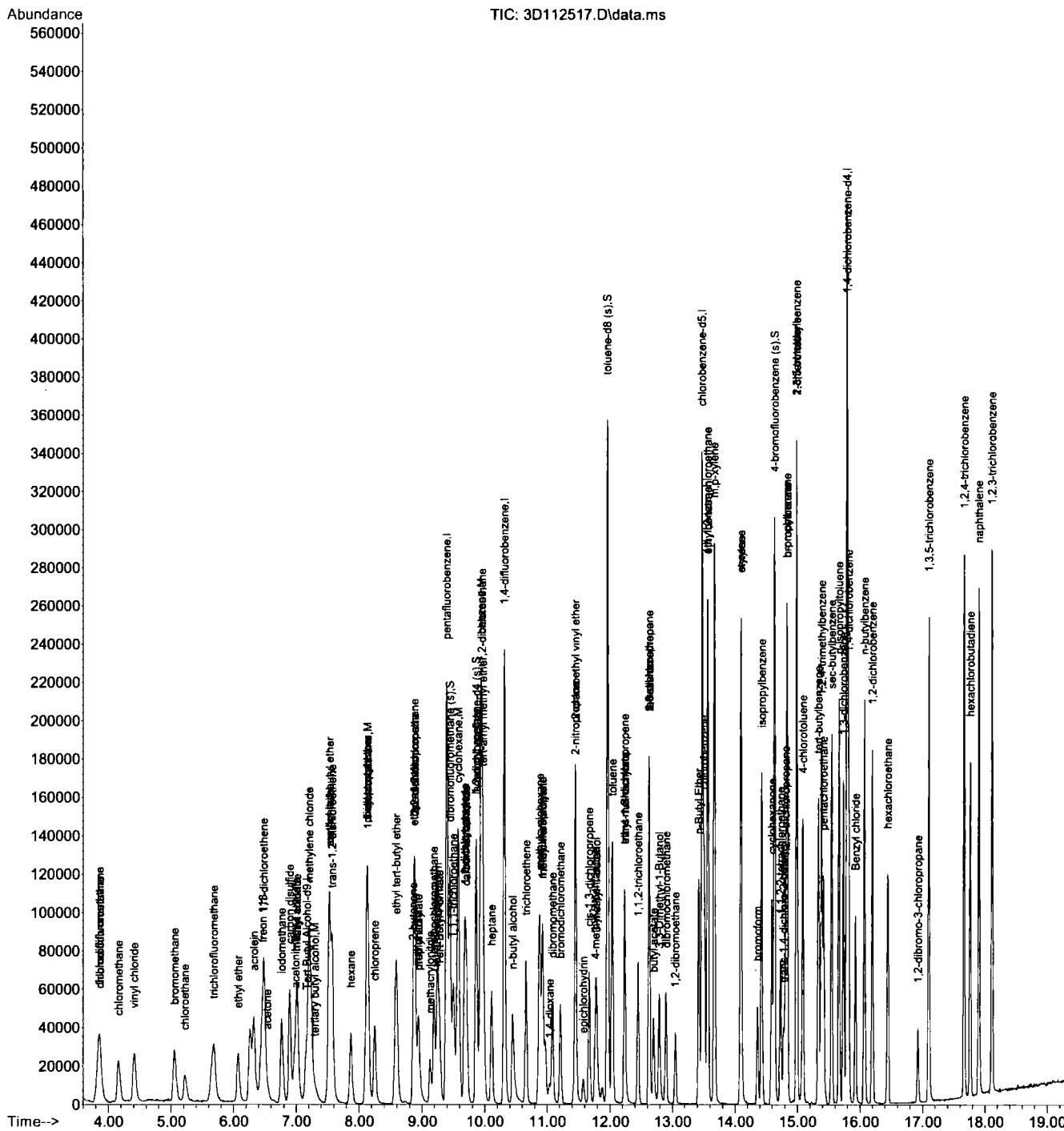
(#) = qualifier out of range (m) = manual integration (+) = signals summed

7.6.12
7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
Data File : 3D112517.D
Acq On : 11 Aug 2015 11:16 am
Operator : ximenac
Sample : cc4810-20
Misc : MS89341,V3D4822,5,,,1
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 11 16:10:05 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112542.D
 Acq On : 11 Aug 2015 11:17 pm
 Operator : ximenac
 Sample : cc4810-50
 Misc : MS89468,V3D4823,5,,,1
 ALS Vial : 35 Sample Multiplier: 1

Quant Time: Aug 12 09:32:24 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) Tert Butyl Alcohol-d9	7.173	65	90136	500.00	ug/L	0.00
4) pentafluorobenzene	9.391	168	190390	50.00	ug/L	0.00
58) 1,4-difluorobenzene	10.308	114	217047	50.00	ug/L	0.00
88) chlorobenzene-d5	13.475	117	183008	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.787	152	113415	50.00	ug/L	0.00
System Monitoring Compounds						
50) dibromofluoromethane (s)	9.443	113	79309	54.37	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	108.74%	
51) 1,2-dichloroethane-d4 (s)	9.863	65	73799	54.61	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	109.22%	
80) toluene-d8 (s)	11.960	98	246124	51.78	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	103.56%	
106) 4-bromofluorobenzene (s)	14.629	95	88049	50.46	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	100.92%	
Target Compounds						
				Qvalue		
2) 1,4-dioxane	11.037	88	21610	1483.76	ug/L	96
3) tertiary butyl alcohol	7.283	59	49609	254.84	ug/L	88
8) chlorodifluoromethane	3.849	51	148076	57.29	ug/L	99
9) dichlorodifluoromethane	3.838	85	199307	57.60	ug/L	99
10) chloromethane	4.164	50	152786	55.53	ug/L	98
11) vinyl chloride	4.415	62	175548	63.61	ug/L	99
12) bromomethane	5.044	94	104409	58.56	ug/L	97
13) chloroethane	5.212	64	68070	58.23	ug/L	98
16) trichlorofluoromethane	5.674	101	189373	65.76	ug/L	100
18) ethyl ether	6.067	74	37557	52.80	ug/L	98
21) acrolein	6.318	56	142341	498.48	ug/L	98
22) 1,1-dichloroethene	6.481	61	130742	57.73	ug/L	96
23) acetone	6.539	43	23143	54.25	ug/L	93
24) allyl chloride	7.011	76	39285	55.85	ug/L	93
25) acetonitrile	6.969	40	80333	566.31	ug/L	85
27) iodomethane	6.764	142	193438	60.07	ug/L	97
28) iso-butyl alcohol	9.684	41	24991	527.25	ug/L	92
29) carbon disulfide	6.890	76	342752	59.85	ug/L	99
30) methylene chloride	7.199	84	97758	57.61	ug/L	97
31) methyl acetate	6.995	43	52884	55.80	ug/L	97
32) methyl tert butyl ether	7.530	73	274088	54.52	ug/L	98
33) trans-1,2-dichloroethene	7.572	61	118788	57.49	ug/L	98
34) di-isopropyl ether	8.122	45	269440	50.70	ug/L	96
35) ethyl tert-butyl ether	8.583	59	271003	54.13	ug/L	99
36) 2-butanone	8.846	72	7227	53.23	ug/L	94
37) 1,1-dichloroethane	8.138	63	143209	58.34	ug/L	99
38) chloroprene	8.248	53	94040	52.34	ug/L	96
39) acrylonitrile	7.524	53	142489	290.01	ug/L	98
40) vinyl acetate	8.133	86	10225	56.06	ug/L	97
41) ethyl acetate	8.861	45	8499	45.51	ug/L #	22
42) 2,2-dichloropropane	8.877	77	138578	53.94	ug/L	97
43) cis-1,2-dichloroethene	8.877	96	86974	52.42	ug/L	99
44) propionitrile	8.935	54	103282	564.11	ug/L	92
45) methyl acrylate	8.940	55	68126	55.20	ug/L	98
46) bromochloromethane	9.186	128	46883	58.71	ug/L	95
47) tetrahydrofuran	9.228	42	23521	48.91	ug/L	99
48) chloroform	9.244	83	135486	57.24	ug/L	97
49) Tert-Butyl Formate	9.270	59	72249	52.43	ug/L #	97
52) freon 113	6.449	151	86161	61.80	ug/L	99
53) methacrylonitrile	9.118	41	35698	49.91	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112542.D
 Acq On : 11 Aug 2015 11:17 pm
 Operator : ximenac
 Sample : cc4810-50
 Misc : MS89468,V3D4823,5,,,1
 ALS Vial : 35 Sample Multiplier: 1

Quant Time: Aug 12 09:32:24 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
54) 1,1,1-trichloroethane	9.496	97	149830	61.31	ug/L	99
57) tert-amyl methyl ether	9.968	73	266691	53.29	ug/L	96
59) epichlorohydrin	11.572	57	29781	249.94	ug/L	97
60) n-butyl alcohol	10.439	56	93525	2639.10	ug/L	96
61) cyclohexane	9.569	84	143558	56.48	ug/L	95
62) carbon tetrachloride	9.705	117	138066	65.77	ug/L	99
63) 1,1-dichloropropene	9.674	75	91709	58.66	ug/L	98
64) hexane	7.870	57	68892	49.34	ug/L	97
65) benzene	9.936	78	287622	56.40	ug/L	99
66) heptane	10.109	57	43439	54.76	ug/L	96
67) isopropyl acetate	9.857	43	116582	52.55	ug/L	99
68) 1,2-dichloroethane	9.952	62	88641	58.19	ug/L	97
69) trichloroethene	10.660	95	69910	56.70	ug/L	98
72) 2-nitropropane	11.436	41	23097	51.24	ug/L	97
73) 2-chloroethyl vinyl ether	11.446	63	213337	272.53	ug/L	99
74) methyl methacrylate	10.922	41	93021	53.35	ug/L	96
75) 1,2-dichloropropane	10.917	63	74932	56.21	ug/L	95
76) methylcyclohexane	10.875	83	141959	61.54	ug/L	98
77) dibromomethane	11.079	93	48138	61.38	ug/L	95
78) bromodichloromethane	11.210	83	94624	58.33	ug/L	100
79) cis-1,3-dichloropropene	11.672	75	108083	56.76	ug/L	96
81) 4-methyl-2-pentanone	11.761	58	25813	53.19	ug/L	98
82) toluene	12.033	92	164755	58.79	ug/L	99
83) 3-methyl-1-butanol	11.782	70	37649	1026.69	ug/L	97
84) trans-1,3-dichloropropene	12.233	75	99241	58.64	ug/L	97
85) ethyl methacrylate	12.227	69	73497	49.14	ug/L	99
86) 1,1,2-trichloroethane	12.448	83	53544	56.87	ug/L	96
87) 2-hexanone	12.621	58	18518	46.68	ug/L	96
89) tetrachloroethene	12.621	166	82914	53.21	ug/L	97
90) 1,3-dichloropropane	12.626	76	97175	54.73	ug/L	98
91) butyl acetate	12.694	56	37786	51.42	ug/L	97
92) 3,3-Dimethyl-1-Butanol	12.788	57	74574	424.57	ug/L	96
93) dibromochloromethane	12.893	129	82021	59.23	ug/L	99
94) 1,2-dibromoethane	13.045	107	64974	55.64	ug/L	99
95) n-Butyl Ether	13.423	57	263624	47.62	ug/L	98
96) chlorobenzene	13.507	112	185944	56.66	ug/L	98
97) 1,1,1,2-tetrachloroethane	13.564	131	87180	59.90	ug/L	99
98) ethylbenzene	13.559	91	304305	54.13	ug/L	99
99) m,p-xylene	13.664	106	236694	112.59	ug/L	99
100) o-xylene	14.083	106	121982	56.81	ug/L	98
101) styrene	14.094	104	190505	52.73	ug/L	99
103) bromoform	14.361	173	61700	57.11	ug/L	98
105) isopropylbenzene	14.424	105	326905	57.44	ug/L	99
107) bromobenzene	14.823	156	92605	55.13	ug/L	96
108) cyclohexanone	14.587	55	43290	257.85	ug/L	96
109) 1,1,2,2-tetrachloroethane	14.728	83	92291	56.03	ug/L	96
110) trans-1,4-dichloro-2-b...	14.770	53	15239	41.80	ug/L	96
111) 1,2,3-trichloropropane	14.802	110	21896	58.21	ug/L	90
112) n-propylbenzene	14.833	91	354174	57.16	ug/L	100
114) 2-chlorotoluene	14.980	126	82628	57.08	ug/L	98
115) 4-chlorotoluene	15.080	126	79561	56.61	ug/L	94
116) 1,3,5-trimethylbenzene	14.985	105	284456	58.73	ug/L	98
117) tert-butylbenzene	15.331	119	232172	58.18	ug/L	100
118) pentachloroethane	15.415	167	71757	65.93	ug/L	98
119) 1,2,4-trimethylbenzene	15.378	105	277948	57.16	ug/L	97
120) sec-butylbenzene	15.546	105	383709	60.17	ug/L	98
121) 1,3-dichlorobenzene	15.735	146	178922	57.28	ug/L	98
122) p-isopropyltoluene	15.662	119	322743	57.96	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
 Data File : 3D112542.D
 Acq On : 11 Aug 2015 11:17 pm
 Operator : ximenac
 Sample : cc4810-50
 Misc : MS89468,V3D4823,5,,,1
 ALS Vial : 35 Sample Multiplier: 1

Quant Time: Aug 12 09:32:24 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev (Min)
123) 1,4-dichlorobenzene	15.814	146	182339	54.42	ug/L	100
124) 1,2-dichlorobenzene	16.191	146	192134	57.06	ug/L	98
126) n-butylbenzene	16.065	92	163765	56.20	ug/L	99
128) 1,2-dibromo-3-chloropr...	16.920	157	25408	51.63	ug/L	93
129) 1,3,5-trichlorobenzene	17.093	180	205876	57.59	ug/L	100
130) 1,2,4-trichlorobenzene	17.659	180	213462	58.09	ug/L	97
131) hexachlorobutadiene	17.759	225	103316	56.80	ug/L	99
132) naphthalene	17.900	128	443378	56.96	ug/L	100
133) 1,2,3-trichlorobenzene	18.110	180	224592	58.86	ug/L	99
134) hexachloroethane	16.443	119	69405	60.34	ug/L	98
135) Benzyl chloride	15.924	91	149724	45.76	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

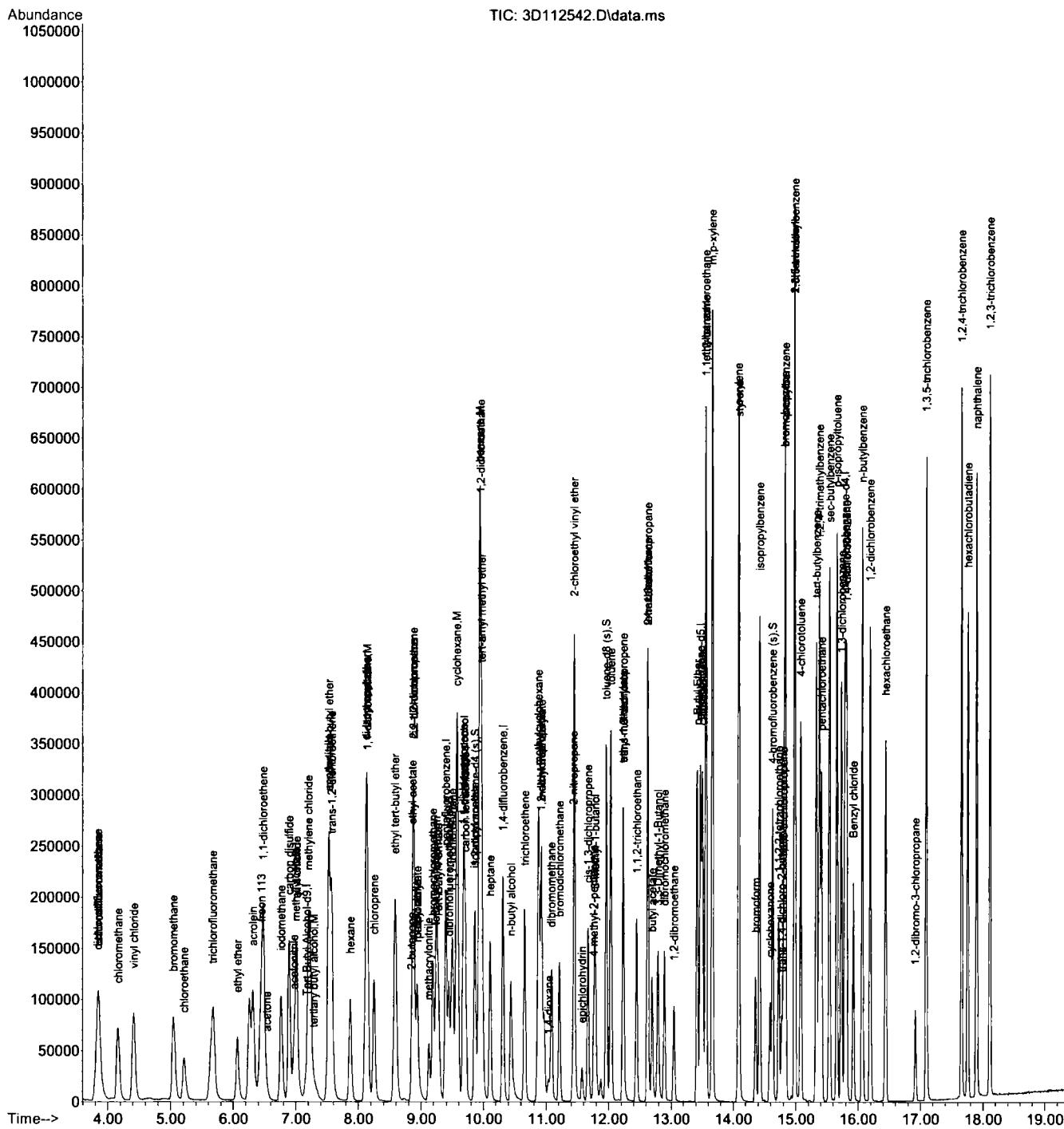
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Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\3D\v3d4822\
Data File : 3D112542.D
Acq On : 11 Aug 2015 11:17 pm
Operator : ximenac
Sample : cc4810-50
Misc : MS89468,V3D4823,5,,,1
ALS Vial : 35 Sample Multiplier: 1

Quant Time: Aug 12 09:32:24 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112563.D
 Acq On : 12 Aug 2015 9:17 am
 Operator : ximenac
 Sample : cc4810-20
 Misc : MS89457,V3D4824,5,,,1
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 12 14:29:03 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Tert Butyl Alcohol-d9	7.173	65	97817	500.00	ug/L	0.00
4) pentafluorobenzene	9.401	168	204794	50.00	ug/L	0.01
58) 1,4-difluorobenzene	10.319	114	227986	50.00	ug/L	0.01
88) chlorobenzene-d5	13.480	117	187452	50.00	ug/L	0.00
104) 1,4-dichlorobenzene-d4	15.793	152	110302	50.00	ug/L	0.00

System Monitoring Compounds

50) dibromofluoromethane (s)	9.449	113	82653	52.68	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 105.36%	
51) 1,2-dichloroethane-d4 (s)	9.868	65	76752	52.80	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 105.60%	
80) toluene-d8 (s)	11.965	98	256659	51.40	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 102.80%	
106) 4-bromofluorobenzene (s)	14.634	95	86012	50.68	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 101.36%	

Target Compounds

					Qvalue	
2) 1,4-dioxane	11.042	88	9959	630.10	ug/L	91
3) tertiary butyl alcohol	7.288	59	20296	96.07	ug/L	80
8) chlorodifluoromethane	3.859	51	57601	20.72	ug/L	97
9) dichlorodifluoromethane	3.854	85	82562	22.18	ug/L	97
10) chloromethane	4.169	50	59192	20.00	ug/L	99
11) vinyl chloride	4.410	62	65708	22.14	ug/L	98
12) bromomethane	5.060	94	43160	22.50	ug/L	98
13) chloroethane	5.228	64	27840	22.14	ug/L	98
16) trichlorofluoromethane	5.684	101	75619	24.41	ug/L	98
18) ethyl ether	6.072	74	15913	20.80	ug/L	92
21) acrolein	6.329	56	56011	182.35	ug/L	97
22) 1,1-dichloroethene	6.492	61	48634	19.97	ug/L	95
23) acetone	6.549	43	10066	21.94	ug/L	88
24) allyl chloride	7.021	76	15306	20.23	ug/L	91
25) acetonitrile	6.979	40	31641	207.37	ug/L	83
27) iodomethane	6.769	142	71808	20.73	ug/L	96
28) iso-butyl alcohol	9.695	41	10383	203.65	ug/L	97
29) carbon disulfide	6.895	76	129102	20.96	ug/L	100
30) methylene chloride	7.210	84	37608	20.60	ug/L	95
31) methyl acetate	7.005	43	21579	21.17	ug/L	97
32) methyl tert butyl ether	7.540	73	100994	18.68	ug/L	100
33) trans-1,2-dichloroethene	7.582	61	46266	20.82	ug/L	97
34) di-isopropyl ether	8.127	45	99826	17.46	ug/L	93
35) ethyl tert-butyl ether	8.594	59	97865	18.17	ug/L	99
36) 2-butanone	8.856	72	2825	19.35	ug/L	86
37) 1,1-dichloroethane	8.148	63	57519	21.78	ug/L	96
38) chloroprene	8.258	53	35047	18.13	ug/L	96
39) acrylonitrile	7.535	53	58132	110.00	ug/L	90
40) vinyl acetate	8.138	86	4127	21.04	ug/L	98
41) ethyl acetate	8.872	45	3263	16.24	ug/L	# 1
42) 2,2-dichloropropane	8.882	77	60173	21.77	ug/L	96
43) cis-1,2-dichloroethene	8.888	96	33920	19.01	ug/L	97
44) propionitrile	8.945	54	42657	216.60	ug/L	85
45) methyl acrylate	8.950	55	27367	20.61	ug/L	98
46) bromochloromethane	9.192	128	18615	21.67	ug/L	96
47) tetrahydrofuran	9.234	42	9682	18.72	ug/L	97
48) chloroform	9.249	83	53449	20.99	ug/L	98
49) Tert-Butyl Formate	9.281	59	26203	17.68	ug/L	# 95
52) freon 113	6.465	151	32611	21.74	ug/L	99
53) methacrylonitrile	9.134	41	14362	18.67	ug/L	87
54) 1,1,1-trichloroethane	9.506	97	59177	22.51	ug/L	97

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112563.D
 Acq On : 12 Aug 2015 9:17 am
 Operator : ximenac
 Sample : cc4810-20
 Misc : MS89457,V3D4824,5,,,1
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 12 14:29:03 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
57) tert-amyl methyl ether	9.973	73	95408	17.72	ug/L	96
59) epichlorohydrin	11.582	57	12448	99.46	ug/L	97
60) n-butyl alcohol	10.455	56	38286	1028.52	ug/L	99
61) cyclohexane	9.574	84	57490	21.53	ug/L	94
62) carbon tetrachloride	9.705	117	53940	24.46	ug/L	97
63) 1,1-dichloropropene	9.679	75	36284	22.10	ug/L	97
64) hexane	7.881	57	25914	17.67	ug/L	97
65) benzene	9.941	78	114725	21.42	ug/L	99
66) heptane	10.114	57	15993	19.20	ug/L	94
67) isopropyl acetate	9.868	43	46031	19.75	ug/L	95
68) 1,2-dichloroethane	9.957	62	34988	21.87	ug/L	97
69) trichloroethene	10.665	95	27784	21.45	ug/L	99
72) 2-nitropropane	11.446	41	9339	19.72	ug/L	89
73) 2-chloroethyl vinyl ether	11.457	63	82696	100.57	ug/L	96
74) methyl methacrylate	10.927	41	38862	21.22	ug/L	96
75) 1,2-dichloropropane	10.927	63	30503	21.78	ug/L	90
76) methylcyclohexane	10.885	83	51374	21.20	ug/L	99
77) dibromomethane	11.084	93	19062	23.14	ug/L	85
78) bromodichloromethane	11.215	83	37030	21.73	ug/L	97
79) cis-1,3-dichloropropene	11.677	75	41112	20.55	ug/L	97
81) 4-methyl-2-pentanone	11.766	58	10236	20.08	ug/L	96
82) toluene	12.044	92	64614	21.95	ug/L	96
83) 3-methyl-1-butanol	11.792	70	14297	371.17	ug/L	96
84) trans-1,3-dichloropropene	12.238	75	38377	21.59	ug/L	98
85) ethyl methacrylate	12.238	69	26603	16.93	ug/L	96
86) 1,1,2-trichloroethane	12.453	83	21341	21.58	ug/L	95
87) 2-hexanone	12.631	58	7127	17.10	ug/L	97
89) tetrachloroethene	12.631	166	32054	20.08	ug/L	98
90) 1,3-dichloropropane	12.636	76	37877	20.83	ug/L	96
91) butyl acetate	12.699	56	13449	17.87	ug/L	89
92) 3,3-Dimethyl-1-Butanol	12.794	57	26532	147.47	ug/L	97
93) dibromochloromethane	12.898	129	31470	22.19	ug/L	98
94) 1,2-dibromoethane	13.051	107	25700	21.49	ug/L	96
95) n-Butyl Ether	13.428	57	93964	16.57	ug/L	96
96) chlorobenzene	13.512	112	72931	21.69	ug/L	96
97) 1,1,1,2-tetrachloroethane	13.570	131	33378	22.39	ug/L	97
98) ethylbenzene	13.564	91	116180	20.18	ug/L	99
99) m,p-xylene	13.674	106	90653	42.10	ug/L	99
100) o-xylene	14.089	106	45211	20.56	ug/L	94
101) styrene	14.104	104	68449	18.50	ug/L	96
103) bromoform	14.361	173	23452	21.19	ug/L	99
105) isopropylbenzene	14.429	105	117275	21.19	ug/L	99
107) bromobenzene	14.833	156	35580	21.78	ug/L	93
108) cyclohexanone	14.597	55	12069	113.68	ug/L	97
109) 1,1,2,2-tetrachloroethane	14.734	83	36431	22.74	ug/L	99
110) trans-1,4-dichloro-2-b...	14.781	53	5908	16.66	ug/L	84
111) 1,2,3-trichloropropane	14.807	110	8445	23.09	ug/L	86
112) n-propylbenzene	14.838	91	130317	21.62	ug/L	99
114) 2-chlorotoluene	14.985	126	30970	22.00	ug/L	93
115) 4-chlorotoluene	15.085	126	29434	21.53	ug/L	91
116) 1,3,5-trimethylbenzene	14.990	105	103124	21.89	ug/L	98
117) tert-butylbenzene	15.336	119	80629	20.77	ug/L	98
118) pentachloroethane	15.415	167	26490	25.03	ug/L	94
119) 1,2,4-trimethylbenzene	15.384	105	100133	21.17	ug/L	98
120) sec-butylbenzene	15.551	105	136274	21.97	ug/L	98
121) 1,3-dichlorobenzene	15.735	146	66388	21.85	ug/L	95
122) p-isopropyltoluene	15.667	119	115732	21.37	ug/L	100
123) 1,4-dichlorobenzene	15.819	146	67677	20.77	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
 Data File : 3D112563.D
 Acq On : 12 Aug 2015 9:17 am
 Operator : ximenac
 Sample : cc4810-20
 Misc : MS89457,V3D4824,5,,,1
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 12 14:29:03 2015
 Quant Method : C:\msdchem\1\METHODS\M3D4810.M
 Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
 QLast Update : Thu Aug 06 11:10:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
124) 1,2-dichlorobenzene	16.196	146	70488	21.52	ug/L	97
126) n-butylbenzene	16.071	92	57533	20.30	ug/L	96
128) 1,2-dibromo-3-chloropr...	16.925	157	9161	19.14	ug/L	93
129) 1,3,5-trichlorobenzene	17.098	180	73721	21.20	ug/L	99
130) 1,2,4-trichlorobenzene	17.664	180	75868	21.23	ug/L	95
131) hexachlorobutadiene	17.764	225	37716	21.32	ug/L	96
132) naphthalene	17.900	128	158867	20.98	ug/L	100
133) 1,2,3-trichlorobenzene	18.115	180	82043	22.11	ug/L	99
134) hexachloroethane	16.443	119	24527	21.93	ug/L	98
135) Benzyl chloride	15.929	91	60346	18.97	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

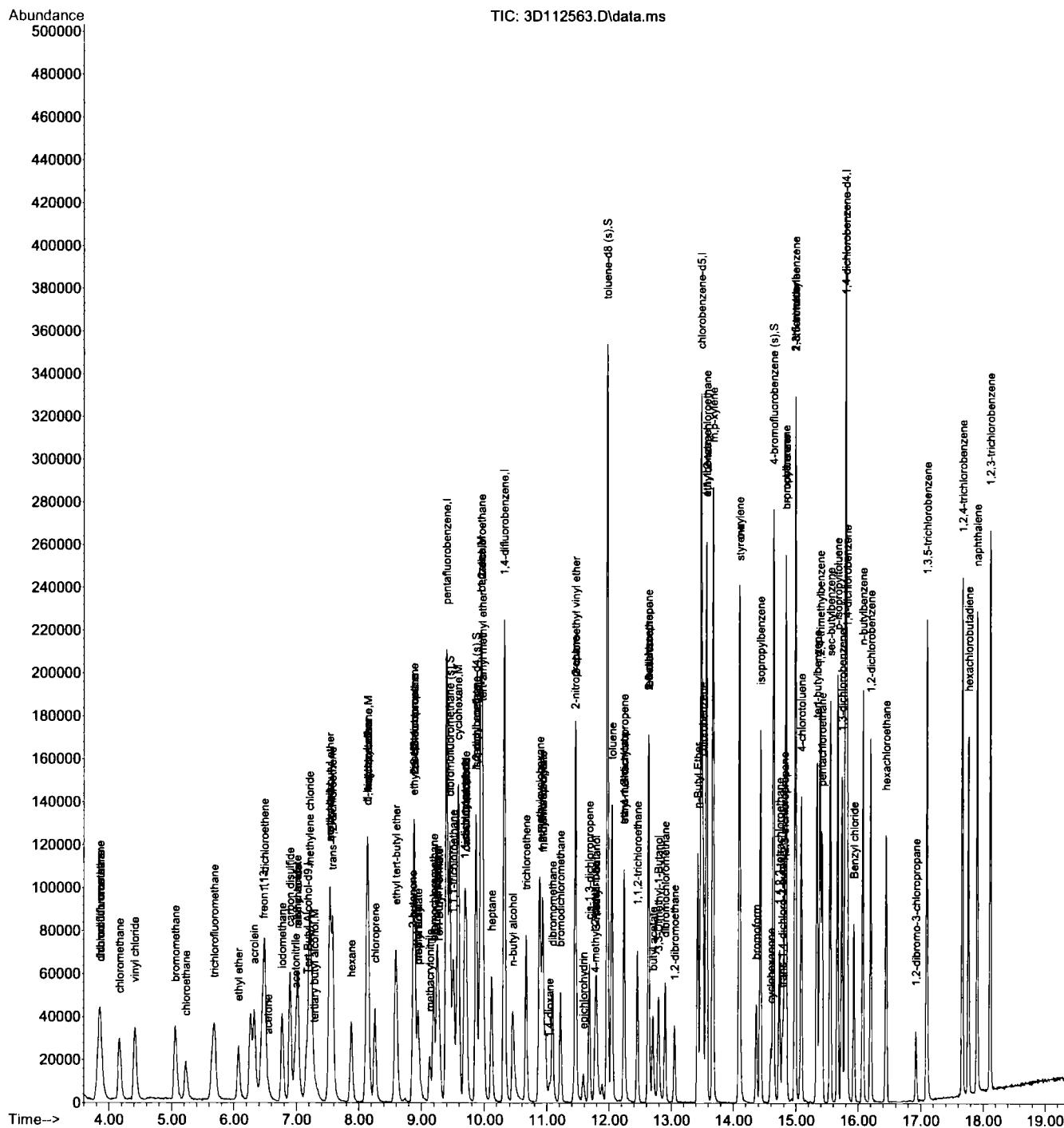
7614

7

Quantitation Report (QT Reviewed)

Data Path : C:\msdchem\1\DATA\
Data File : 3D112563.D
Acq On : 12 Aug 2015 9:17 am
Operator : ximenac
Sample : cc4810-20
Misc : MS89457,V3D4824,5,,,1
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 12 14:29:03 2015
Quant Method : C:\msdchem\1\METHODS\M3D4810.M
Quant Title : Method SW846 8260B, ZB 624 60m x 0.25mm x 1.4 um
QLast Update : Thu Aug 06 11:10:54 2015
Response via : Initial Calibration





ACCU TEST

VOLATILE ANALYSIS LOG

Batch ID: V3D4810

Print Analyst Name: KIMENA

Analyst Signature: Greg Holladay

Columns: ~~2B-624~~ (60 m x 0.25 mm i.d. 4 µm)

Method V826c

Initial Cal. Method M3D4810

Standards

Standard Data

Lot #	Description	Conc.
W015-2114	-41.79	STD A
W015-2114	-72.16	STD B
W015-2114	-71.11	STD C
W015-2114	-54.9	STD A/R
W015-2114	-63	IIS

Standard Data			
Lot #	Description	Conc.	
W015-2114	49.20	Ext A	100 ppm
W015-2114	73.15	Ext B	100 ppm
W015-2114	70.6	Ext C	100 ppm
W015-2114	48.4	Ext A2	1000 ppm
W015-2114	51.4	HOK	100 ppm

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature: _____

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Date: 7/20/15

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt. (ul)	Secondary dilution	L + I S U	S U	Status (Date)	Comments	pH* <2
	3D112234	bfb										OK		
	112235	IC4810-0.2	V8260C A2					5				W	OK 1 ml ABC Acr → some	
	112236	IC4810-0.5		✓				5				W	OK 2.5 ml ABC Acr → some	
	112237	IC4810-1		✓				5				W	OK 200 ml Acr + 2 ml ABC → 200 ml	
	112238	IC4810-2		✓				5				W	OK 200 ml Acr + 2 ml ABC → some	
	112239	IC4810-5		✓				5				W	OK 5 ml ABC Acr → some	
	112240	IC4810-10		✓				5				W	OK 5 ml ABC Acr → some	
	112241	IC4810-20		✓				5				W	OK 10 ml ABC Acr → some	
	112242	IC4810-50		✓				5				W	OK 25 ml ABC Acr → some	
	112243	IC4810-100		✓				5				W	OK 1 ml Acr + 50 ml ABC → some	
	112244	IC4810-200		✓				5				W	OK 100 ml ABC → some	
	112245	ib						5						
	112246	ib						5						
	112247	ICV4810-50		✓				5				W	OK 25 ml Ext (ABC Acr H2O) → some	
													KC 7/19/15	

MTX = Matrix Designate W for water, S for soil, O for oil **L+** = Library Search **IS = Internal Standard Area** **SU = Surrogate**

Sample Amt = Volume (ML) or Weight (g) MOH amt = volume (mL) extract injected. IF pH > 2 comment on sample result

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error.

Form: QB001-9

Rev. Date: 2/14/2007

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VOLATILE ANALYSIS LOG

Date: 8/11/15

Batch ID: V3D4822

Print Analyst Name: XIMENA

Analyst Signature: Guy Hollis

Columns: ZB-624 (60m x 0.25mm x 1.4um)

Method V826OC

Initial Cal. Method H3D4810

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature: TM Date: 8/11/15

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt (ul)	Secondary dilution	L + I S U	S U	Status (Data)	Comments	pH* <2
	3D117S15	bfb										OK	9:53 AM	
	112516	CC4810-20					5					NG	Compounds ↑	
	112517	CC4810-20					5					OK	20mL Std ABC Acet /100ml 10↓ 23↑	
	112518	ib					5					W	OK	
	112519	mb					5					W	OK	
	112520	bs					5					W	OK	25mL EST(ABC Acet hex) /100ml
	112521	ib					5					W	OK	
	112522	JC1106-7	89468	SL	1		5		IX			W	OK	✓
	112523	JC1106-5		G	1		5		IX			W	OK	✓
	112524	JC1106-13		G	1		5		IX			W	OK	✓
	112525	JC1106-10		G	1		5		IX			W	OK	✓
	112526	JC1106-13 ms		G	6		5		IX			W	OK	DEVAL ABC Acet
	112527	JC1106-13 msd		G	7		5		IX			W	OK	to 50 mL
	112528	ib					5					W		
	112529	JC1106-19	89468	SL	1		5		IX			W	OK	✓
	112530	JC1106-6		G	1		5		IX			W	OK	✓
	112531	JC1106-3		G	1		5		IX			W	OK	✓

MTX = Matrix Designate W for water, S for soil, O for oil. L+ =Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt.= volume (ul) extract injected * IF pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9
Rev. Date: 2/14/2007

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ACCUTEST.

VOLATILE ANALYSIS LOG

Date: 8/11/15

Standard Data		
Lot #	Description	Cone.

Standard Data		
Lot #	Description	Cone.

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature: JM Date: 8/10/15

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt. (uL)	Secondary dilution	L I S U	Status (Data)	Comments	pH* < 2
	3D112532	JC 1106 - 4	89468 SL	G J	1		5		IX		W/OE		✓
	112 533	JC 1106 - 15	✓	G J	1		5		IX		W/OE		✓
	112 534	JC 1106 - 18	✓	G J	1		5		IX		W/OE		✓
	112 535	JC 1106 - 8	✓	G J	2		5		IX		W/OE		✓
	112 536	JC 1106 - 14	✓	G J	1		5		IX		W/OE		✓
	112 537	JC 1106 - 12	✓	G J	1		5		IX		W/OE		✓
	112 538	JC 1106 - 11	✓	G J	1		5		IX		W/OE		✓
	112 539	JC 1106 - 9	✓	G J	2		5		IX		W/OE out of detection: 55 ppm		✓
	112 540	ib					5						
XLC 8/11/15													

MTX = Matrix Designate W for water, S for soil, O for oil. L+ =Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt= volume (uL) extract injected * IF pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9
Rev. Date: 2/14/2007

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ACCUTEST.

VOLATILE ANALYSIS LOG

Date: 8/12/15

Standard Data

Lot #	Description	Conc.
V01T-2114-95-11	GAT A	100 ppm
V01T-2114-72-64	GAT B	100 ppm
V01T-2114-111-2	GAT C	100 ppm
V01T-2114-107-4	GAT A&B	1000 ppm
V01T-2114-92-2	Hex	100 ppm

Standard Data

Lot #	Description	Conc.
V01T-2114-93-6	STD A	100 ppm
V01T-2114-72-46	STD B	100 ppm
V01T-2114-114-2	STD C	100 ppm
V01T-2114-54-41	STD A&B	1000 ppm
V01T-2114-89	TIS	250/500 ppm

Batch ID: V3D4823 / V3D4824

Print Analyst Name: Ximeena

Analyst Signature: Mylottab

Columns: 78-624 (60 mm x 25 mm x 1.4 μm)

Method U8260C

Initial Cal. Method M3D4810

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044. PH paper 212713

Supervisor Signature: Sup S/11/16

HHS

Date: 8/12/15

7.7.3
7

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (mL or g)	MOH amt. (μL)	Secondary dilution	L + S U	I S U	Status (Date)	Comments	pH* <2
	3D 112562	bfb								O			8:16 am	
	112563	-CLU810-20					5			W	O	C	2, 16, 62 ↑ 20 mL Std A B C A&B 100%	
	112564	ib					5			W	O	C		
	112565	mb2					5			W	O	C		
	112566	JC 1072-1 ms	89457 TELnot TBA	G	4		5		IX	W	O	C	STWABCABD	✓
	112567*	JC 1072-1 msd	✓	G	4		5		IX	W	O	C	630 mL	✓
	112568	ib					5			W				
	112569	mb					5			W	O	C		
	112570	bs					5			W	O	C	25 mL Ext (A B C A&B) 150 mL	
	112571	ib					5			W				
R	112572	JC 1106-9	89468 SL	G	1		5		IX	W	O	C		✓
R	112573	JC 1106 -1b	✓	G	2		5		IX	W	O	C		✓
R	112574	JC 1106 -2	✓	G	2		5		IX	W	O	C		✓
R	112575	JC 1106 -1	✓	G	2		5		IX	W	O	C		✓
R	112576	JC 726-8	89228 TELnot TBA	G	2		5		IX	W	O	C		5
R	112577	JC 1082-3	89462 TELnot TBA	G	1		10/50		5X	W	O	C	/DL 50X	✓
R	112578	JC 726-9	89228 TELnot TBA	G	2		5		IX	W	O	C		✓
R	112579	JC 726-2	✓	G	2		5		IX	W	O	C		✓

MTX = Matrix Designate W for water, S for soil, O for oil. L+ = Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (mL) or Weight (g); MOH amt = volume (μL) extract injected * IF pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9
Rev. Date: 2/14/2007

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ACCUTEST.

VOLATILE ANALYSIS LOG

Date: 8/12/15

Standard Data

Lot #	Description	Conc.
VOLT-21M	GAT-A	100 ppm
VOLT-21M	GAT-B	100 ppm
VOLT-21M	GAT-C	100 ppm
VOLT-21M	GAT-AIR	1000 ppm
VOLT-21M	Hex	100 ppm

Standard Data

Lot #	Description	Conc.
VOLT-21M	93.6	STD A
VOLT-21M	72.46	STD B
VOLT-21M	118.2	STD C
VOLT-21M	54.41	STD AIR
VOLT-21M	89	IIS

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044. PH paper 21273

Supervisor Signature: *[Signature]*

Run:

Date: 8/12/15

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt. (ul)	Secondary dilution	L I S + S U	Status (Data)	Comments	pH* <2
	3D 112562	bfb									OK	8:16 am	
	112563	-CLU4810-20					5				W OK	2, 16, 62 ↑ 20 ml Std A B C Acet 100ml	
	112564	ib					5				W OK		
	112565	mb2					5				W OK		
	112566	JC 1072-1 ms	89467 TEL20+TBA	G ✓	4	4	5		IX		W OK	✓	
	112567	JC 1072-1 msd	✓	G	4	4	5		IX		W OK	650ML.	✓
	112568	ib					5				W		
	112569	mb					5				W OK		
	112570	bs					5				W OK	25 ml Ext (A B C Acet) 100ml	
	112571	ib					5				W		
R	112572	JC 1106-9	89468 TEL20+TBA	G ✓	1	1	5		IX		W OK		✓
R	112573	JC 1106-16	✓	G	2	2	5		IX		W OK		✓
R	112574	JC 1106-2	✓	G	2	2	5		IX		W OK		✓
R	112575	JC 1106-1	✓	G	2	2	5		IX		W OK		✓
R	112576	JC 726-8	89228 TEL20+TBA	G ✓	2	2	5		IX		W OK		5
R	112577	JC 1082-3	89462 TEL20+TBA	G ✓	1	1	10 / 50		5X		W OK	/DL 50X	✓
R	112578	JC 726-9	89228 TEL20+TBA	G ✓	2	2	5		IX		W OK		✓
R	112579	JC 726-2	✓	G	2	5	5		IX		W OK		

MTX = Matrix Designate W for water, S for soil, O for oil. L+ = Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt = volume (ul) extract injected * IF pH > 2, comment on sample result.

- All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9
Rev. Date: 2/14/2007

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VOLATILE ANALYSIS LOG

Date: 8/12/05

Batch ID: V3D4824

Standard Data

Lot #	Description	Conc.

Standard Data

Lot #	Description	Conc.

See 8/12/05

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature: HWS Date: 8/13/05

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt. (ul)	Secondary dilution	L + I S U	Status (Data)	Comments	pH* <2
	3D112580	JC1106-9	SL	G	2		5		IX	W	OC		✓
	112581	JC1106-9	gsc	G	2		5		IX	W	OC		✓
DY	112582	JC11082	89462 TU20+TBA	G	1		1/50		50X	W	OC		✓
	112583	JC11082-3	89517 PTLL42+TBA	G	2		5		IX	W	OC		✓
	112584	JC1177-1	HDXYL,NAP OXYL,TBA	E	F		5		IX	W	OC		5
	112585	JC1182-2		E	N	7	5		IX	W	RR	IRRIK 40	✓
	112586	JC1182-1		E	F	6	5		IX	W	PR	IRRIK 40 7:46pm	✓
	112587	bfb										OC	
	112588	CC4810-50					5			W	OC	20ml Std A+B C/Aw/150ml 16↑ 52,62,76,133↑	
	112589	CC4810-50					5			W	OC		
	112590	mb2					5			W	OC		
	112591	JC1006-2	89463 TU20+TBA	G	1		5		IX	W	OC		✓
	112592	JC1006-1	TBA	G	2		5		IX	W	OC		✓
	112593	JC1080-2	89461 TU20+TBA	G	1		5		IX	W	OC		✓
	112594	JC1080-1		G	1		5		IX	W	OC		✓
	112595	JC1082-2	89462 TU20+	G	1		5		IX	W	OC		✓
	112596	JC1082-1	TBA	G	1		5		IX	W	OC		✓
	112597	JC1080-3	89461 TU20+TBA	G	1		5		IX	W	OC		✓

MTX = Matrix Designate W for water, S for soil, O for oil. L+ =Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt.= volume (ul) extract injected * IF pH > 2, comment on sample result.

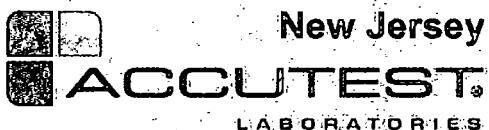
All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculations; 4 = analyst's correction error

HWS 8/13/05

Form: OR001-9

Rev. Date: 2/14/2007

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08/24/15

Technical Report for

United Technologies Corporation

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

60339110 PO#59048ACM

Accutest Job Number: JC1107

Sampling Dates: 08/06/15 - 08/07/15

Report to:

**AECOM, INC.
27755 Diehl Road Suite 100
Warrenville, IL 60555
peter.hollatz@aecom.com**

ATTN: Peter Hollatz

Total number of pages in report: 215



Test results contained within this data package meet the requirements of the National Environmental Laboratory Accreditation Program and/or state specific certification programs as applicable.

Nancy T. Cole

Nancy Cole
Laboratory Director

Client Service contact: Marie Meidhof 732-329-0200

Certifications: NJ(12129), NY(10983), CA, CT, DE, FL, IL, IN, KS, KY, LA, MA, MD, MI, MT, NC, OH VAP (CL0056), AK (UST-103), AZ (AZ0786), PA, RI, SC, TN, TX, VA, WV, DoD ELAP (L-A-B L2248)

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Test results relate only to samples analyzed.

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Sample Summary

United Technologies Corporation

Job No: JC1107

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Collected Date	Time By	Matrix Received	Code Type	Client Sample ID	
JC1107-1	08/06/15	12:10 AH	08/08/15	AQ	Ground Water	HSSER-RAMW01-080615
JC1107-2	08/06/15	11:05 AH	08/08/15	AQ	Field Blank Water	HSSER-FBLK02-080615
JC1107-3	08/06/15	13:20 AH	08/08/15	AQ	Ground Water	HSSER-RAMW02-080615
JC1107-3D	08/06/15	13:20 AH	08/08/15	AQ	Water Dup/MSD	HSSER-MSD02-080615
JC1107-3S	08/06/15	13:20 AH	08/08/15	AQ	Water Matrix Spike	HSSER-MS02-080615
JC1107-4	08/06/15	14:50 AH	08/08/15	AQ	Ground Water	HSSER-RAMW03-080615
JC1107-5	08/06/15	16:10 AH	08/08/15	AQ	Ground Water	HSSER-RAMW08-080615
JC1107-6	08/06/15	00:00 AH	08/08/15	AQ	Ground Water	HSSER-DUP02-080615
JC1107-7	08/07/15	08:10 AH	08/08/15	AQ	Ground Water	HSSER-RAMW07-080715
JC1107-8	08/07/15	09:35 AH	08/08/15	AQ	Ground Water	HSSER-RAMW06-080715
JC1107-9	08/07/15	10:55 AH	08/08/15	AQ	Ground Water	HSSER-RAMW05-080715
JC1107-10	08/07/15	11:55 AH	08/08/15	AQ	Ground Water	HSSER-RAMW04-080715
JC1107-11	08/07/15	11:15 AH	08/08/15	AQ	Equipment Blank	HSSER-EBLK02-080715

Sample Summary

(continued)

United Technologies Corporation

Job No: JC1107

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Collected Date	Time By	Matrix Received	Code Type	Client Sample ID
JC1107-12	08/07/15	11:15 AH	08/08/15	AQ	Trip Blank Water HSSER-TRIP02-080415



CASE NARRATIVE / CONFORMANCE SUMMARY

Client: United Technologies Corporation

Job No JC1107

Site: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Report Date 8/24/2015 11:09:50 A

On 08/08/2015, 9 Sample(s), 1 Trip Blank(s) and 1 Field Blank(s) and 1 Equipment Blank(s) were received at Accutest Laboratories at a maximum corrected temperature of 1.9 C. Samples were intact and chemically preserved, unless noted below. An Accutest Job Number of JC1107 was assigned to the project. Laboratory sample ID, client sample ID and dates of sample collection are detailed in the report's Results Summary Section.

Specified quality control criteria were achieved for this job except as noted below. For more information, please refer to the analytical results and QC summary pages.

Volatiles by GCMS By Method SW846 8260C

Matrix: AQ

Batch ID: V4B2296

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JC1107-3MS, JC1107-3MSD were used as the QC samples indicated.

Matrix: AQ

Batch ID: V4B2297

- All samples were analyzed within the recommended method holding time.
- All method blanks for this batch meet method specific criteria.
- Sample(s) JC869-3DUP, JC869-4MS were used as the QC samples indicated.

Accutest certifies that data reported for samples received, listed on the associated custody chain or analytical task order, were produced to specifications meeting Accutest's Quality System precision, accuracy and completeness objectives except as noted.

Estimated non-standard method measurement uncertainty data is available on request, based on quality control bias and implicit for standard methods. Acceptable uncertainty requires tested parameter quality control data to meet method criteria.

Accutest Laboratories is not responsible for data quality assumptions if partial reports are used and recommends that this report be used in its entirety. Data release is authorized by Accutest Laboratories indicated via signature on the report cover

Summary of Hits

Page 1 of 2

Job Number: JC1107

Account: United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Collected: 08/06/15 thru 08/07/15



Lab Sample ID Analyte	Client Sample ID	Result/ Qual	RL	MDL	Units	Method
JC1107-1 HSSER-RAMW01-080615						
1,1-Dichloroethane	0.0037	0.0010	0.00017	mg/l	SW846 8260C	
cis-1,2-Dichloroethene	0.00070 J	0.0010	0.00027	mg/l	SW846 8260C	
Tetrachloroethene	0.0052	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0039	0.0010	0.00025	mg/l	SW846 8260C	
1,1,2-Trichloroethane	0.00022 J	0.0010	0.00021	mg/l	SW846 8260C	
Trichloroethene	0.0012	0.0010	0.00022	mg/l	SW846 8260C	
JC1107-2 HSSER-FBLK02-080615						
Toluene	0.00033 J	0.0010	0.00016	mg/l	SW846 8260C	
JC1107-3 HSSER-RAMW02-080615						
1,1-Dichloroethane	0.0029	0.0010	0.00017	mg/l	SW846 8260C	
Tetrachloroethene	0.0025	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0032	0.0010	0.00025	mg/l	SW846 8260C	
Trichloroethene	0.00041 J	0.0010	0.00022	mg/l	SW846 8260C	
JC1107-4 HSSER-RAMW03-080615						
1,1-Dichloroethane	0.00039 J	0.0010	0.00017	mg/l	SW846 8260C	
Tetrachloroethene	0.0011	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.00043 J	0.0010	0.00025	mg/l	SW846 8260C	
Trichloroethene	0.00043 J	0.0010	0.00022	mg/l	SW846 8260C	
JC1107-5 HSSER-RAMW08-080615						
1,1-Dichloroethane	0.00025 J	0.0010	0.00017	mg/l	SW846 8260C	
JC1107-6 HSSER-DUP02-080615						
1,1-Dichloroethane	0.00041 J	0.0010	0.00017	mg/l	SW846 8260C	
Tetrachloroethene	0.0011	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.00041 J	0.0010	0.00025	mg/l	SW846 8260C	
Trichloroethene	0.00044 J	0.0010	0.00022	mg/l	SW846 8260C	
JC1107-7 HSSER-RAMW07-080715						
1,1-Dichloroethane	0.0072	0.0010	0.00017	mg/l	SW846 8260C	
1,1-Dichloroethene	0.0069	0.0010	0.00051	mg/l	SW846 8260C	
cis-1,2-Dichloroethene	0.0016	0.0010	0.00027	mg/l	SW846 8260C	
Tetrachloroethene	0.0015	0.0010	0.00040	mg/l	SW846 8260C	
1,1,1-Trichloroethane	0.0762	0.0010	0.00025	mg/l	SW846 8260C	

Summary of Hits

Job Number: JC1107

Account: United Technologies Corporation

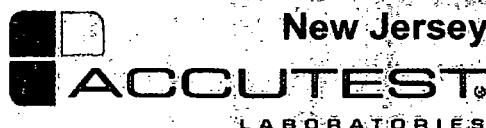
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Collected: 08/06/15 thru 08/07/15



Lab Sample ID Analyte	Client Sample ID	Result/ Qual	RL	MDL	Units	Method
Trichloroethene		0.0015	0.0010	0.00022	mg/l	SW846 8260C
JC1107-8 HSSER-RAMW06-080715						
1,1-Dichloroethane		0.0026	0.0010	0.00017	mg/l	SW846 8260C
1,1-Dichloroethene		0.0056	0.0010	0.00051	mg/l	SW846 8260C
cis-1,2-Dichloroethene		0.0068	0.0010	0.00027	mg/l	SW846 8260C
1,1,1-Trichloroethane		0.0596	0.0010	0.00025	mg/l	SW846 8260C
Trichloroethene		0.00061 J	0.0010	0.00022	mg/l	SW846 8260C
JC1107-9 HSSER-RAMW05-080715						
1,1,1-Trichloroethane		0.00052 J	0.0010	0.00025	mg/l	SW846 8260C
JC1107-10 HSSER-RAMW04-080715						
Tetrachloroethene		0.00043 J	0.0010	0.00040	mg/l	SW846 8260C
1,1,1-Trichloroethane		0.00039 J	0.0010	0.00025	mg/l	SW846 8260C
JC1107-11 HSSER-EBLK02-080715						
Toluene		0.00032 J	0.0010	0.00016	mg/l	SW846 8260C
JC1107-12 HSSER-TRIP02-080415						

No hits reported in this sample.



Sample Results

Report of Analysis

Accutest Laboratories

Report of Analysis

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Client Sample ID:	HSSER-RAMW01-080615	Date Sampled:	08/06/15
Lab Sample ID:	JC1107-1	Date Received:	08/08/15
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Method:	SW846 8260C		
Project:	ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54425.D	1	08/12/15	TP	n/a	n/a	V4B2296
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0037	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.00070	0.0010	0.00027	mg/l	J
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0052	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0039	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	0.00022	0.0010	0.00021	mg/l	J
79-01-6	Trichloroethene	0.0012	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	97%		76-120%
17060-07-0	1,2-Dichloroethane-D4	92%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	105%		78-117%

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-FBLK02-080615**Lab Sample ID:** JC1107-2**Date Sampled:** 08/06/15**Matrix:** AQ - Field Blank Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54426.D	1	08/12/15	TP	n/a	n/a	V4B2296
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	0.00033	0.0010	0.00016	mg/l	J
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	97%		76-120%
17060-07-0	1,2-Dichloroethane-D4	94%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	104%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-RAMW02-080615
Lab Sample ID: JC1107-3
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/06/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54416.D	1	08/12/15	TP	n/a	n/a	V4B2296
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0029	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0025	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0032	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.00041	0.0010	0.00022	mg/l	J
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	98%		76-120%
17060-07-0	1,2-Dichloroethane-D4	92%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	104%		78-117%

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-RAMW03-080615**Lab Sample ID:** JC1107-4**Date Sampled:** 08/06/15**Matrix:** AQ - Ground Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54427.D	1	08/12/15	TP	n/a	n/a	V4B2296
Run #2							

Purge Volume

Run #1 5.0 ml

Run #2

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.00039	0.0010	0.00017	mg/l	J
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0011	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.00043	0.0010	0.00025	mg/l	J
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.00043	0.0010	0.00022	mg/l	J
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	97%		76-120%
17060-07-0	1,2-Dichloroethane-D4	93%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	104%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-RAMW08-080615
Lab Sample ID: JC1107-5
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/06/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54428.D	1	08/12/15	TP	n/a	n/a	V4B2296
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.00025	0.0010	0.00017	mg/l	J
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	98%		76-120%
17060-07-0	1,2-Dichloroethane-D4	94%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	104%		78-117%

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID:	HSSER-DUP02-080615	Date Sampled:	08/06/15
Lab Sample ID:	JC1107-6	Date Received:	08/08/15
Matrix:	AQ - Ground Water	Percent Solids:	n/a
Method:	SW846 8260C		
Project:	ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL		

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54429.D	1	08/12/15	TP	n/a	n/a	V4B2296
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.00041	0.0010	0.00017	mg/l	J
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0011	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.00041	0.0010	0.00025	mg/l	J
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.00044	0.0010	0.00022	mg/l	J
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	98%		76-120%
17060-07-0	1,2-Dichloroethane-D4	94%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	103%		78-117%

ND = Not detected MDL = Method Detection Limit
 RL = Reporting Limit
 E = Indicates value exceeds calibration range

J = Indicates an estimated value
 B = Indicates analyte found in associated method blank
 N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

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Client Sample ID: HSSER-RAMW07-080715
Lab Sample ID: JC1107-7
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/07/15
Date Received: 08/08/15
Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54451.D	1	08/13/15	TP	n/a	n/a	V4B2297
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0072	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	0.0069	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.0016	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.0015	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0762	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.0015	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	97%		76-120%
17060-07-0	1,2-Dichloroethane-D4	92%		73-122%
2037-26-5	Toluene-D8	98%		84-119%
460-00-4	4-Bromofluorobenzene	104%		78-117%

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: HSSER-RAMW06-080715**Lab Sample ID:** JC1107-8**Date Sampled:** 08/07/15**Matrix:** AQ - Ground Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54450.D	1	08/13/15	TP	n/a	n/a	V4B2297
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	0.0026	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	0.0056	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	0.0068	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.0596	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	0.00061	0.0010	0.00022	mg/l	J
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	97%		76-120%
17060-07-0	1,2-Dichloroethane-D4	91%		73-122%
2037-26-5	Toluene-D8	98%		84-119%
460-00-4	4-Bromofluorobenzene	107%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: HSSER-RAMW05-080715
Lab Sample ID: JC1107-9
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/07/15
Date Received: 08/08/15
Percent Solids: n/a

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54452.D	1	08/13/15	TP	n/a	n/a	V4B2297
Run #2							

Run #	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.00052	0.0010	0.00025	mg/l	J
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	97%		76-120%
17060-07-0	1,2-Dichloroethane-D4	93%		73-122%
2037-26-5	Toluene-D8	98%		84-119%
460-00-4	4-Bromofluorobenzene	106%		78-117%

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: HSSER-RAMW04-080715
Lab Sample ID: JC1107-10
Matrix: AQ - Ground Water
Method: SW846 8260C
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/07/15
Date Received: 08/08/15
Percent Solids: n/a

Run #	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54453.D	1	08/13/15	TP	n/a	n/a	V4B2297
Run #2							

Purge Volume	
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	0.00043	0.0010	0.00040	mg/l	J
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	0.00039	0.0010	0.00025	mg/l	J
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	98%		76-120%
17060-07-0	1,2-Dichloroethane-D4	94%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	105%		78-117%

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: HSSER-EBLK02-080715

Lab Sample ID: JC1107-11

Matrix: AQ - Equipment Blank

Method: SW846 8260C

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Date Sampled: 08/07/15

Date Received: 08/08/15

Percent Solids: n/a

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54423.D	1	08/12/15	TP	n/a	n/a	V4B2296
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	0.00032	0.0010	0.00016	mg/l	J
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	96%		76-120%
17060-07-0	1,2-Dichloroethane-D4	91%		73-122%
2037-26-5	Toluene-D8	100%		84-119%
460-00-4	4-Bromofluorobenzene	104%		78-117%

ND = Not detected MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

Accutest Laboratories

Report of Analysis

Page 1 of 1

Client Sample ID: HSSER-TRIP02-080415**Lab Sample ID:** JC1107-12**Date Sampled:** 08/07/15**Matrix:** AQ - Trip Blank Water**Date Received:** 08/08/15**Method:** SW846 8260C**Percent Solids:** n/a**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
Run #1	4B54424.D	1	08/12/15	TP	n/a	n/a	V4B2296
Run #2							

	Purge Volume
Run #1	5.0 ml
Run #2	

VOA Special List

CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	0.0010	0.00017	mg/l	
107-06-2	1,2-Dichloroethane	ND	0.0010	0.00018	mg/l	
75-35-4	1,1-Dichloroethene	ND	0.0010	0.00051	mg/l	
156-59-2	cis-1,2-Dichloroethene	ND	0.0010	0.00027	mg/l	
156-60-5	trans-1,2-Dichloroethene	ND	0.0010	0.00065	mg/l	
100-41-4	Ethylbenzene	ND	0.0010	0.00027	mg/l	
75-09-2	Methylene chloride	ND	0.0020	0.00073	mg/l	
127-18-4	Tetrachloroethene	ND	0.0010	0.00040	mg/l	
108-88-3	Toluene	ND	0.0010	0.00016	mg/l	
71-55-6	1,1,1-Trichloroethane	ND	0.0010	0.00025	mg/l	
79-00-5	1,1,2-Trichloroethane	ND	0.0010	0.00021	mg/l	
79-01-6	Trichloroethene	ND	0.0010	0.00022	mg/l	
75-01-4	Vinyl chloride	ND	0.0010	0.00015	mg/l	

CAS No.	Surrogate Recoveries	Run# 1	Run# 2	Limits
1868-53-7	Dibromofluoromethane	96%		76-120%
17060-07-0	1,2-Dichloroethane-D4	92%		73-122%
2037-26-5	Toluene-D8	99%		84-119%
460-00-4	4-Bromofluorobenzene	105%		78-117%

ND = Not detected

MDL = Method Detection Limit

J = Indicates an estimated value

RL = Reporting Limit

B = Indicates analyte found in associated method blank

E = Indicates value exceeds calibration range

N = Indicates presumptive evidence of a compound

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Misc. Forms



Custody Documents and Other Forms

Includes the following where applicable:

- Chain of Custody
- Sample Tracking Chronicle
- Internal Chain of Custody



GW
WTB

CHAIN OF CUSTODY

PAGE 1 OF 2

2235 Route 130, Dayton, NJ 08810
TEL: 732-329-0200 FAX: 732-329-3499/3480
www.accutest.com

FED-Ex Tracking #	Box/Carton #
1634940616410	Accutest Job # JC1107

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)		Matrix Codes					
Company Name ACOM	Project Name UTAS Plants V2 Feb 17 yes.	Street Address 27755 D.oh1 Rd Suite 100	City Kennebunkville IL	State IL	Zip 60555						
Street Address 27755 D.oh1 Rd Suite 100	City Kentwood	State IL	Billing Information (if different from Report to)		Company Name						
Project Contact Peter Hollatz Peter.Hollatz@acom.com	Project # 600339110	Phone # (630)836-5380 (630)836-1711	City 69048 ACOM	State IL	Zip 60555	Street Address					
Sampler(s) Name(s) Peter Hollatz	Phone # (630)836-1760	Project Manager Peter Hollatz	City 69048 ACOM		State IL	Zip 60555	Attention:				
Across Sample #	Field ID / Point of Collection	MEOHDI Val #	Collection				Number of preserved Bottles				
-1	HSSER-RAMW01-080615	8-6-15	1210	AH GW	3	X		X	VOC	LAB USE ONLY	V859
-2	HSSER-FALK02-080615	8-6-15	1105		3	X		X			
-3	HSSER-RA1102-080615	8-6-15	1320		3	X		X			
-3	HSSER-MSD02-080615	8-6-15	1320		3	y		X			
	HSSER-MSD02-080615	8-6-15	1320		3	y		X			
-4	HSSER-RAMW03-080615	8-6-15	1450		3	X		X			
-5	HSSER-RAMW08-080615	8-6-15	1610		3	X		X			
-6	HSSER-DUP02-080615	8-6-15	0000		3	X		X		INITIAL ASSESSMENT 3A DOM	
-7	HSSER-RAMW07-080715	8-7-15	0810		3	X		X		LABEL VERIFICATION JMW	
-8	HSSER-RAMW06-080715	8-7-15	0935		3	X		X			
-9	HSSER-RAMW05-080715	8-7-15	1055		3	X		X			
-10	HSSER-RAMW04-080715	8-7-15	1155		3	X		X			
Turnaround Time (Business days)		Data Deliverable Information		Comments / Special Instructions							
<input type="checkbox"/> Std. 10 Business Days <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day RUSH <input type="checkbox"/> 2 Day RUSH <input type="checkbox"/> 1 Day RUSH <input type="checkbox"/> other _____		<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULL/TI (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NY Data of Known Quality Protocol Reporting		<input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input type="checkbox"/> Other _____							
Emergency & Rush T/A data available via LabLink		Commercial "A" = Results Only, Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data		Comments / Special Instructions							
Sample Custody must be documented below each time samples change possession, including courier delivery.											
Ratiquested by Sampler: 1 Peter Hollatz	Date Time: 8-7-15 1800	Received By: 1 FX	Ratiquested By: 2	Received By: FX	Date Time: 8-8-15 10:00	Received By: 2					
Ratiquested by Sampler: 3	Date Time: 	Received By: 3	Ratiquested By: 4	Received By: 	Date Time: 8-8-15 10:00	Received By: 4					
Ratiquested by: 5	Date Time: 	Received By: 5	Custody Seal #	<input type="checkbox"/> intact <input type="checkbox"/> Not intact	Preserved where applicable	On Ice <input checked="" type="checkbox"/>	Cooler Temp 32.2 + ip				

JC1107: Chain of Custody

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CHAIN OF CUSTODY

PAGE 2 OF 2

2235 Route 130, Dayton, NJ 08810
TEL. 732-329-0200 FAX. 732-329-3499/3480
www.acutesol.com

Client / Reporting Information		Project Information		Requested Analysis (see TEST CODE sheet)		Matrix Codes										
Company Name AECOM	Project Name: UTAS Plants 1/2 Facil:1/4r					DW - Drinking Water GW - Ground Water WW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rouse Blank TB - Trip Blank										
Street Address 27755 Ditch Rd. Suite 100	Sample	Billing Information (if different from Report to)														
City Lakeville IL	State IL	City Portfield	State FL	Company Name												
Project Contact Peter Hollatz	E-mail Peter.Hollatz@AECOM.com	Project # 10339110	Street Address													
Phone # 630-836-5786	Fax # 630-836-1711	Client Purchased Order # 59048 ACAN	City	State	Zip											
Sample(s) Name(s) Allen Hollatz	Phone # 630-836-1700	Project Manager Peter Hollatz	Attention:													
Acquisition Sample #		Field ID / Point of Collection	Collection	Number of preserved bottles												
		METHOD Var #	Date	Sampled by	Metric	# of bottles	HQ	HC	HR	NC	ME	GW	MECH	ENONE	LOC	LAB USE ONLY
-11		HSSER-EBLK02-080715	8-7-15	1115	AH	GW	3	X							X	V859
-12		HSSFR-Tip02-080415	8-4-15	—			2	X						X		
Turnaround Time (Business days)		Approved By (Accruent Pd): Date:	Data Deliverable Information		Comments / Special Instructions											
<input type="checkbox"/> Std. 10 Business Days <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day RUSH <input type="checkbox"/> 2 Day RUSH <input type="checkbox"/> 1 Day RUSH <input type="checkbox"/> other _____		<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLY! (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input type="checkbox"/> Other _____ <input type="checkbox"/> NJ Date of Known Quality Protocol Reporting Commercial "A" = Results Only, Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	* List of 13 UCC Level IV Data													
Emergency & Rush T/A data available VIA Lablink		Sample Custody must be documented below each time samples change possession, including courier delivery.														
Retainership by Sampler: 1 CECIL HOLLATZ	Date Time: 8-7-15 1800	Received By: 1 FX	Retainership By: 2	Date Time: 8-7-15 10:00	Received By: 2											
Retainership by Sampler: 3	Date Time: 8-7-15 1800	Received By: 3	Retainership By: 4	Date Time: 8-7-15 1800	Received By: 4											
Retainership by Sampler: 5	Date Time: 8-7-15 1800	Received By: 5	Custody Seal #	Preserved where applicable		On Ice: <input checked="" type="checkbox"/>	Carrier Temp: 22C									

JC1107: Chain of Custody
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Accutest Laboratories Sample Receipt Summary

Accutest Job Number: JC1107 Client: _____ Project: _____

Date / Time Received: 8/8/2015 10:00:00 AM Delivery Method: _____ Airbill #: _____

Cooler Temps (Raw Measured) °C: Cooler 1: (2.2);

Cooler Temps (Corrected) °C: Cooler 1: (1.9);

Cooler Security		Y or N		Y or N		Sample Integrity - Documentation		Y or N	
1. Custody Seals Present:	<input checked="" type="checkbox"/> <input type="checkbox"/>	3. COC Present:	<input checked="" type="checkbox"/> <input type="checkbox"/>	1. Sample labels present on bottles:	<input checked="" type="checkbox"/> <input type="checkbox"/>	2. Container labeling complete:	<input checked="" type="checkbox"/> <input type="checkbox"/>	3. Sample container label / COC agree:	<input checked="" type="checkbox"/> <input type="checkbox"/>
2. Custody Seals Intact:	<input checked="" type="checkbox"/> <input type="checkbox"/>	4. Smpl Dates/Time OK	<input checked="" type="checkbox"/> <input type="checkbox"/>						
Cooler Temperature		Y or N				Sample Integrity - Condition		Y or N	
1. Temp criteria achieved:	<input checked="" type="checkbox"/> <input type="checkbox"/>					1. Sample recvd within HT:	<input checked="" type="checkbox"/> <input type="checkbox"/>	2. All containers accounted for:	<input checked="" type="checkbox"/> <input type="checkbox"/>
2. Cooler temp verification:		IR Gun				3. Condition of sample:		Intact	
3. Cooler media:		Ice (Bag)							
4. No. Coolers:		1							
Quality Control Preservation		Y or N		N/A		Sample Integrity - Instructions		Y or N	
1. Trip Blank present / cooler:	<input checked="" type="checkbox"/> <input type="checkbox"/>					1. Analysis requested is clear:	<input checked="" type="checkbox"/> <input type="checkbox"/>	2. Bottles received for unspecified tests	<input type="checkbox"/> <input checked="" type="checkbox"/>
2. Trip Blank listed on COC:	<input checked="" type="checkbox"/> <input type="checkbox"/>					3. Sufficient volume recvd for analysis:	<input checked="" type="checkbox"/> <input type="checkbox"/>	4. Compositing instructions clear:	<input type="checkbox"/> <input type="checkbox"/> <input checked="" type="checkbox"/>
3. Samples preserved properly:	<input checked="" type="checkbox"/> <input type="checkbox"/>					5. Filtering instructions clear:	<input type="checkbox"/> <input type="checkbox"/>		<input checked="" type="checkbox"/>
4. VOCs headspace free:	<input checked="" type="checkbox"/> <input type="checkbox"/>								

Comments

Accutest Laboratories
V:732.329.0200

2235 US Highway 130
F: 732.329.3499

Dayton, New Jersey
www.accutest.com

JC1107: Chain of Custody

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Internal Sample Tracking Chronicle

United Technologies Corporation

Job No: JC1107

ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Method	Analyzed	By	Prepped	By	Test Codes
JC1107-1	Collected: 06-AUG-15 12:10 By: AH HSSER-RAMW01-080615			Received: 08-AUG-15 By: DG		
JC1107-1	SW846 8260C	12-AUG-15 18:46	TP			V8260SL
JC1107-2	Collected: 06-AUG-15 11:05 By: AH HSSER-FBLK02-080615			Received: 08-AUG-15 By: DG		
JC1107-2	SW846 8260C	12-AUG-15 19:14	TP			V8260SL
JC1107-3	Collected: 06-AUG-15 13:20 By: AH HSSER-RAMW02-080615			Received: 08-AUG-15 By: DG		
JC1107-3	SW846 8260C	12-AUG-15 14:35	TP			V8260SL
JC1107-4	Collected: 06-AUG-15 14:50 By: AH HSSER-RAMW03-080615			Received: 08-AUG-15 By: DG		
JC1107-4	SW846 8260C	12-AUG-15 19:42	TP			V8260SL
JC1107-5	Collected: 06-AUG-15 16:10 By: AH HSSER-RAMW08-080615			Received: 08-AUG-15 By: DG		
JC1107-5	SW846 8260C	12-AUG-15 20:10	TP			V8260SL
JC1107-6	Collected: 06-AUG-15 00:00 By: AH HSSER-DUP02-080615			Received: 08-AUG-15 By: DG		
JC1107-6	SW846 8260C	12-AUG-15 20:38	TP			V8260SL
JC1107-7	Collected: 07-AUG-15 08:10 By: AH HSSER-RAMW07-080715			Received: 08-AUG-15 By: DG		
JC1107-7	SW846 8260C	13-AUG-15 06:51	TP			V8260SL
JC1107-8	Collected: 07-AUG-15 09:35 By: AH HSSER-RAMW06-080715			Received: 08-AUG-15 By: DG		
JC1107-8	SW846 8260C	13-AUG-15 06:23	TP			V8260SL

Internal Sample Tracking Chronicle

United Technologies Corporation

Job No: JC1107

ENSRLW: UTAS Plants 1/2 Facility, Rockford, IL
Project No: 60339110 PO#59048ACM

Sample Number	Method	Analyzed	By	Prepped	By	Test Codes
JC1107-9	Collected: 07-AUG-15 10:55 By: AH HSSER-RAMW05-080715			Received: 08-AUG-15 By: DG		
JC1107-9	SW846 8260C	13-AUG-15 07:19	TP			V8260SL
JC1107-10	Collected: 07-AUG-15 11:55 By: AH HSSER-RAMW04-080715			Received: 08-AUG-15 By: DG		
JC1107-10	SW846 8260C	13-AUG-15 07:48	TP			V8260SL
JC1107-11	Collected: 07-AUG-15 11:15 By: AH HSSER-EBLK02-080715			Received: 08-AUG-15 By: DG		
JC1107-11	SW846 8260C	12-AUG-15 17:51	TP			V8260SL
JC1107-12	Collected: 07-AUG-15 11:15 By: AH HSSER-TRIP02-080415			Received: 08-AUG-15 By: DG		
JC1107-12	SW846 8260C	12-AUG-15 18:18	TP			V8260SL

Accutest Internal Chain of Custody

Page 1 of 2

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Received: 08/08/15

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JC1107-1.1	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-1.1	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-1.1	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-1.1	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-2.1	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-2.1	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-2.1	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-2.1	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-3.3	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-3.3	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-3.3	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-3.3	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-4.1	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-4.1	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-4.1	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-4.1	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-5.2	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-5.2	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-5.2	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-5.2	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-6.2	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-6.2	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-6.2	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-6.2	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-7.2	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-7.2	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-7.2	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-7.2	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-8.2	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-8.2	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-8.2	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-8.2	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-9.2	Secured Storage	Toan Pham	08/12/15 14:09	Retrieve from Storage
JC1107-9.2	Toan Pham	GCMS4B	08/12/15 14:09	Load on Instrument
JC1107-9.2	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-9.2	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage

Accutest Internal Chain of Custody

Page 2 of 2

Job Number: JC1107
Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL
Received: 08/08/15

Sample/Bottle Number	Transfer FROM	Transfer TO	Date/Time	Reason
JC1107-10.2	Secured Storage	Toan Pham	08/12/15 14:09	Retrieve from Storage
JC1107-10.2	Toan Pham	GCMS4B	08/12/15 14:09	Load on Instrument
JC1107-10.2	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-10.2	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-11.2	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-11.2	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-11.2	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-11.2	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage
JC1107-12.2	Secured Storage	Toan Pham	08/12/15 12:13	Retrieve from Storage
JC1107-12.2	Toan Pham	GCMS4B	08/12/15 12:13	Load on Instrument
JC1107-12.2	GCMS4B	Zobia Hamid	08/13/15 09:59	Unload from Instrument
JC1107-12.2	Zobia Hamid	Secured Storage	08/13/15 09:59	Return to Storage

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GC/MS Volatiles



QC Data Summaries

Includes the following where applicable:

- Method Blank Summaries
- Blank Spike Summaries
- Matrix Spike and Duplicate Summaries
- Instrument Performance Checks (BFB)
- Internal Standard Area Summaries
- Surrogate Recovery Summaries
- Initial and Continuing Calibration Summaries

Method Blank Summary**Job Number:** JC1107**Account:** UTC United Technologies Corporation**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V4B2296-MB	4B54409.D	1	08/12/15	TP	n/a	n/a	V4B2296

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC1107-1, JC1107-2, JC1107-3, JC1107-4, JC1107-5, JC1107-6, JC1107-11, JC1107-12

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CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	

CAS No. Surrogate Recoveries Limits

1868-53-7	Dibromofluoromethane	98%	76-120%
17060-07-0	1,2-Dichloroethane-D4	92%	73-122%
2037-26-5	Toluene-D8	98%	84-119%
460-00-4	4-Bromofluorobenzene	105%	78-117%

CAS No. Tentatively Identified Compounds R.T. Est. Conc. Units Q

Total TIC, Volatile 0 ug/l

Method Blank Summary

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V4B2297-MB	4B5443.D	1	08/12/15	TP	n/a	n/a	V4B2297

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC1107-7, JC1107-8, JC1107-9, JC1107-10

6.12



CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	

CAS No.	Surrogate Recoveries	Limits
1868-53-7	Dibromofluoromethane	98%
17060-07-0	1,2-Dichloroethane-D4	93%
2037-26-5	Toluene-D8	98%
460-00-4	4-Bromofluorobenzene	104%

CAS No.	Tentatively Identified Compounds	R.T.	Est. Conc.	Units	Q
	Total TIC, Volatile		0	ug/l	

Method Blank Summary**Job Number:** JC1107**Account:** UTC United Technologies Corporation**Project:** ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V4B2297-MB2	4B54457.D	1	08/13/15	TP	n/a	n/a	V4B2297

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC869-3DUP, JC869-4MS

613



CAS No.	Compound	Result	RL	MDL	Units	Q
75-34-3	1,1-Dichloroethane	ND	1.0	0.17	ug/l	
107-06-2	1,2-Dichloroethane	ND	1.0	0.18	ug/l	
75-35-4	1,1-Dichloroethene	ND	1.0	0.51	ug/l	
156-59-2	cis-1,2-Dichloroethene	ND	1.0	0.27	ug/l	
156-60-5	trans-1,2-Dichloroethene	ND	1.0	0.65	ug/l	
100-41-4	Ethylbenzene	ND	1.0	0.27	ug/l	
75-09-2	Methylene chloride	ND	2.0	0.73	ug/l	
127-18-4	Tetrachloroethene	ND	1.0	0.40	ug/l	
108-88-3	Toluene	ND	1.0	0.16	ug/l	
71-55-6	1,1,1-Trichloroethane	ND	1.0	0.25	ug/l	
79-00-5	1,1,2-Trichloroethane	ND	1.0	0.21	ug/l	
79-01-6	Trichloroethene	ND	1.0	0.22	ug/l	
75-01-4	Vinyl chloride	ND	1.0	0.15	ug/l	

CAS No.	Surrogate Recoveries	Limits
1868-53-7	Dibromofluoromethane	96%
17060-07-0	1,2-Dichloroethane-D4	92%
2037-26-5	Toluene-D8	99%
460-00-4	4-Bromofluorobenzene	106%
		76-120%
		73-122%
		84-119%
		78-117%

Blank Spike Summary

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V4B2296-BS	4B54410.D	1	08/12/15	TP	n/a	n/a	V4B2296

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC1107-1, JC1107-2, JC1107-3, JC1107-4, JC1107-5, JC1107-6, JC1107-11, JC1107-12

6.2.1



CAS No.	Compound	Spike ug/l	BSP ug/l	BSP %	Limits
75-34-3	1,1-Dichloroethane	50	46.7	93	80-125
107-06-2	1,2-Dichloroethane	50	49.7	99	78-131
75-35-4	1,1-Dichloroethene	50	51.1	102	73-127
156-59-2	cis-1,2-Dichloroethene	50	47.9	96	77-118
156-60-5	trans-1,2-Dichloroethene	50	45.4	91	75-118
100-41-4	Ethylbenzene	50	49.9	100	80-118
75-09-2	Methylene chloride	50	49.7	99	75-122
127-18-4	Tetrachloroethene	50	50.6	101	69-138
108-88-3	Toluene	50	48.6	97	80-122
71-55-6	1,1,1-Trichloroethane	50	49.5	99	80-131
79-00-5	1,1,2-Trichloroethane	50	47.9	96	78-122
79-01-6	Trichloroethene	50	50.9	102	83-122
75-01-4	Vinyl chloride	50	43.8	88	57-138

CAS No.	Surrogate Recoveries	BSP	Limits
1868-53-7	Dibromofluoromethane	100%	76-120%
17060-07-0	1,2-Dichloroethane-D4	94%	73-122%
2037-26-5	Toluene-D8	100%	84-119%
460-00-4	4-Bromofluorobenzene	96%	78-117%

* = Outside of Control Limits.

Blank Spike Summary

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
V4B2297-BS	4B54434.D	1	08/12/15	TP	n/a	n/a	V4B2297

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC1107-7, JC1107-8, JC1107-9, JC1107-10

CAS No.	Compound	Spike ug/l	BSP ug/l	BSP %	Limits
75-34-3	1,1-Dichloroethane	50	45.7	91	80-125
107-06-2	1,2-Dichloroethane	50	50.1	100	78-131
75-35-4	1,1-Dichloroethene	50	48.2	96	73-127
156-59-2	cis-1,2-Dichloroethene	50	47.2	94	77-118
156-60-5	trans-1,2-Dichloroethene	50	46.0	92	75-118
100-41-4	Ethylbenzene	50	49.1	98	80-118
75-09-2	Methylene chloride	50	47.9	96	75-122
127-18-4	Tetrachloroethene	50	50.8	102	69-138
108-88-3	Toluene	50	48.7	97	80-122
71-55-6	1,1,1-Trichloroethane	50	52.1	104	80-131
79-00-5	1,1,2-Trichloroethane	50	47.8	96	78-122
79-01-6	Trichloroethene	50	50.4	101	83-122
75-01-4	Vinyl chloride	50	43.1	86	57-138

CAS No.	Surrogate Recoveries	BSP	Limits
1868-53-7	Dibromofluoromethane	100%	76-120%
17060-07-0	1,2-Dichloroethane-D4	95%	73-122%
2037-26-5	Toluene-D8	100%	84-119%
460-00-4	4-Bromofluorobenzene	96%	78-117%

* = Outside of Control Limits.

Matrix Spike Summary

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
JC869-4MS	4B54459.D	1	08/13/15	TP	n/a	n/a	V4B2297
JC869-4	4B54442.D	1	08/13/15	TP	n/a	n/a	V4B2297

The QC reported here applies to the following samples:

Method: SW846 8260C

JC1107-7, JC1107-8, JC1107-9, JC1107-10

6.31



CAS No.	Compound	JC869-4 ug/l	Spike Q ug/l	MS ug/l	MS %	Limits
75-34-3	1,1-Dichloroethane	ND	50	49.0	98	60-129
107-06-2	1,2-Dichloroethane	ND	50	51.9	104	72-133
75-35-4	1,1-Dichloroethene	ND	50	54.5	109	40-137
156-59-2	cis-1,2-Dichloroethene	ND	50	51.9	104	57-128
156-60-5	trans-1,2-Dichloroethene	ND	50	51.8	104	53-128
100-41-4	Ethylbenzene	ND	50	54.6	109	38-139
75-09-2	Methylene chloride	ND	50	50.8	102	63-128
127-18-4	Tetrachloroethene	ND	50	55.0	110	43-145
108-88-3	Toluene	ND	50	53.2	106	51-136
71-55-6	1,1,1-Trichloroethane	ND	50	58.9	118	51-141
79-00-5	1,1,2-Trichloroethane	ND	50	49.1	98	71-127
79-01-6	Trichloroethene	ND	50	55.1	110	55-136
75-01-4	Vinyl chloride	ND	50	54.6	109	34-147

CAS No.	Surrogate Recoveries	MS	JC869-4	Limits
1868-53-7	Dibromofluoromethane	101%	98%	76-120%
17060-07-0	1,2-Dichloroethane-D4	95%	93%	73-122%
2037-26-5	Toluene-D8	99%	99%	84-119%
460-00-4	4-Bromofluorobenzene	97%	105%	78-117%

* = Outside of Control Limits.

Matrix Spike/Matrix Spike Duplicate Summary

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
JC1107-3MS	4B54420.D	1	08/12/15	TP	n/a	n/a	V4B2296
JC1107-3MSD	4B54421.D	1	08/12/15	TP	n/a	n/a	V4B2296
JC1107-3	4B54416.D	1	08/12/15	TP	n/a	n/a	V4B2296

The QC reported here applies to the following samples:

Method: SW846 8260C

JC1107-1, JC1107-2, JC1107-3, JC1107-4, JC1107-5, JC1107-6, JC1107-11, JC1107-12

CAS No.	Compound	JC1107-3		Spike ug/l	MS ug/l	MS %	Spike ug/l	MSD ug/l	MSD %	RPD	Limits
		ug/l	Q								Rec/RPD
75-34-3	1,1-Dichloroethane	2.9	50	50.5	95	50	49.9	94	1	60-129/13	
107-06-2	1,2-Dichloroethane	ND	50	52.3	105	50	51.2	102	2	72-133/12	
75-35-4	1,1-Dichloroethene	ND	50	52.3	105	50	52.1	104	0	40-137/17	
156-59-2	cis-1,2-Dichloroethene	ND	50	50.4	101	50	50.2	100	0	57-128/13	
156-60-5	trans-1,2-Dichloroethene	ND	50	49.7	99	50	49.4	99	1	53-128/15	
100-41-4	Ethylbenzene	ND	50	53.2	106	50	52.5	105	1	38-139/12	
75-09-2	Methylene chloride	ND	50	50.2	100	50	50.0	100	0	63-128/13	
127-18-4	Tetrachloroethene	2.5	50	54.1	103	50	53.1	101	2	43-145/15	
108-88-3	Toluene	ND	50	51.4	103	50	51.0	102	1	51-136/13	
71-55-6	1,1,1-Trichloroethane	3.2	50	58.7	111	50	57.3	108	2	51-141/16	
79-00-5	1,1,2-Trichloroethane	ND	50	50.4	101	50	50.7	101	1	71-127/12	
79-01-6	Trichloroethene	0.41	J	50	53.6	106	50	52.7	105	2	55-136/14
75-01-4	Vinyl chloride	ND	50	42.2	84	50	42.9	86	2	34-147/17	

CAS No.	Surrogate Recoveries	MS	MSD	JC1107-3	Limits
1868-53-7	Dibromofluoromethane	98%	99%	98%	76-120%
17060-07-0	1,2-Dichloroethane-D4	94%	94%	92%	73-122%
2037-26-5	Toluene-D8	98%	99%	99%	84-119%
460-00-4	4-Bromofluorobenzene	95%	95%	104%	78-117%

* = Outside of Control Limits.

Duplicate Summary

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample	File ID	DF	Analyzed	By	Prep Date	Prep Batch	Analytical Batch
JC869-3DUP	4B54458.D	1	08/13/15	TP	n/a	n/a	V4B2297
JC869-3	4B54441.D	1	08/13/15	TP	n/a	n/a	V4B2297

The QC reported here applies to the following samples:**Method: SW846 8260C**

JC1107-7, JC1107-8, JC1107-9, JC1107-10

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CAS No.	Compound	JC869-3		DUP			
		ug/l	Q	ug/l	Q	RPD	Limits
75-34-3	1,1-Dichloroethane	ND		ND		nc	30
107-06-2	1,2-Dichloroethane	ND		ND		nc	30
75-35-4	1,1-Dichloroethene	ND		ND		nc	30
156-59-2	cis-1,2-Dichloroethene	ND		ND		nc	30
156-60-5	trans-1,2-Dichloroethene	ND		ND		nc	30
100-41-4	Ethylbenzene	ND		ND		nc	30
75-09-2	Methylene chloride	ND		ND		nc	30
127-18-4	Tetrachloroethene	ND		ND		nc	30
108-88-3	Toluene	ND		ND		nc	30
71-55-6	1,1,1-Trichloroethane	ND		ND		nc	30
79-00-5	1,1,2-Trichloroethane	ND		ND		nc	30
79-01-6	Trichloroethene	ND		ND		nc	30
75-01-4	Vinyl chloride	ND		ND		nc	30

CAS No.	Surrogate Recoveries	DUP	JC869-3	Limits
1868-53-7	Dibromofluoromethane	99%	98%	76-120%
17060-07-0	1,2-Dichloroethane-D4	92%	93%	73-122%
2037-26-5	Toluene-D8	97%	98%	84-119%
460-00-4	4-Bromofluorobenzene	109%	104%	78-117%

* = Outside of Control Limits.

Instrument Performance Check (BFB)

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample:	V4B2289-BFB	Injection Date:	08/05/15
Lab File ID:	4B54241.D	Injection Time:	09:05
Instrument ID:	GCMS4B		

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	14.99 - 40.0% of mass 95	23821	20.4	Pass
75	30.0 - 60.0% of mass 95	56474	48.5	Pass
95	Base peak, 100% relative abundance	116525	100.0	Pass
96	5.0 - 9.0% of mass 95	7868	6.75	Pass
173	Less than 2.0% of mass 174	0	0.00	(0.00) ^a Pass
174	50.0 - 120.0% of mass 95	116618	100.1	Pass
175	5.0 - 9.0% of mass 174	9749	8.37	(8.36) ^a Pass
176	95.0 - 101.0% of mass 174	112621	96.6	(96.6) ^a Pass
177	5.0 - 9.0% of mass 176	7858	6.74	(6.98) ^b Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V4B2289-IC2289	4B54242.D	08/05/15	09:39	00:34	Initial cal 0.2
V4B2289-IC2289	4B54243.D	08/05/15	10:07	01:02	Initial cal 0.5
V4B2289-IC2289	4B54244.D	08/05/15	10:35	01:30	Initial cal 1
V4B2289-IC2289	4B54245.D	08/05/15	11:03	01:58	Initial cal 2
V4B2289-IC2289	4B54246.D	08/05/15	11:31	02:26	Initial cal 5
V4B2289-IC2289	4B54247.D	08/05/15	11:59	02:54	Initial cal 10
V4B2289-IC2289	4B54248.D	08/05/15	12:27	03:22	Initial cal 20
V4B2289-ICC2289	4B54249.D	08/05/15	12:54	03:49	Initial cal 50
V4B2289-IC2289	4B54250.D	08/05/15	13:22	04:17	Initial cal 100
V4B2289-IC2289	4B54251.D	08/05/15	13:50	04:45	Initial cal 200
V4B2289-ICV2289	4B54254.D	08/05/15	15:20	06:15	Initial cal verification 50

Instrument Performance Check (BFB)

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample: V4B2296-BFB

Injection Date: 08/12/15

Lab File ID: 4B54405.D

Injection Time: 09:13

Instrument ID: GCMS4B

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	14.99 - 40.0% of mass 95	18701	19.3	Pass
75	30.0 - 60.0% of mass 95	44842	46.4	Pass
95	Base peak, 100% relative abundance	96672	100.0	Pass
96	5.0 - 9.0% of mass 95	6517	6.74	Pass
173	Less than 2.0% of mass 174	0	0.00	(0.00) ^a Pass
174	50.0 - 120.0% of mass 95	101530	105.0	Pass
175	5.0 - 9.0% of mass 174	8723	9.02	(8.59) ^a Pass
176	95.0 - 101.0% of mass 174	98101	101.5	(96.6) ^a Pass
177	5.0 - 9.0% of mass 176	6561	6.79	(6.69) ^b Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V4B2296-CC2289	4B54407.D	08/12/15	10:09	00:56	Continuing cal 20
V4B2296-MB	4B54409.D	08/12/15	11:13	02:00	Method Blank
V4B2296-BS	4B54410.D	08/12/15	11:47	02:34	Blank Spike
ZZZZZZ	4B54413.D	08/12/15	13:10	03:57	(unrelated sample)
ZZZZZZ	4B54414.D	08/12/15	13:38	04:25	(unrelated sample)
ZZZZZZ	4B54415.D	08/12/15	14:06	04:53	(unrelated sample)
JC1107-3	4B54416.D	08/12/15	14:35	05:22	HSSER-RAMW02-080615
JC1107-3MS	4B54420.D	08/12/15	16:27	07:14	Matrix Spike
JC1107-3MSD	4B54421.D	08/12/15	16:55	07:42	Matrix Spike Duplicate
JC1107-11	4B54423.D	08/12/15	17:51	08:38	HSSER-EBLK02-080715
JC1107-12	4B54424.D	08/12/15	18:18	09:05	HSSER-TRIP02-080415
JC1107-1	4B54425.D	08/12/15	18:46	09:33	HSSER-RAMW01-080615
JC1107-2	4B54426.D	08/12/15	19:14	10:01	HSSER-FBLK02-080615
JC1107-4	4B54427.D	08/12/15	19:42	10:29	HSSER-RAMW03-080615
JC1107-5	4B54428.D	08/12/15	20:10	10:57	HSSER-RAMW08-080615
JC1107-6	4B54429.D	08/12/15	20:38	11:25	HSSER-DUP02-080615

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Instrument Performance Check (BFB)

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample:	V4B2297-BFB	Injection Date:	08/12/15
Lab File ID:	4B54430.D	Injection Time:	21:06
Instrument ID:	GCMS4B		

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	14.99 - 40.0% of mass 95	18824	19.0	Pass
75	30.0 - 60.0% of mass 95	47162	47.6	Pass
95	Base peak, 100% relative abundance	98989	100.0	Pass
96	5.0 - 9.0% of mass 95	6769	6.84	Pass
173	Less than 2.0% of mass 174	0	0.00	(0.00) ^a Pass
174	50.0 - 120.0% of mass 95	102565	103.6	Pass
175	5.0 - 9.0% of mass 174	8410	8.50	(8.20) ^a Pass
176	95.0 - 101.0% of mass 174	100890	101.9	(98.4) ^a Pass
177	5.0 - 9.0% of mass 176	6717	6.79	(6.66) ^b Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V4B2297-CC2289	4B54431.D	08/12/15	21:35	00:29	Continuing cal 50
V4B2297-MB	4B54433.D	08/12/15	22:30	01:24	Method Blank
V4B2297-BS	4B54434.D	08/12/15	22:58	01:52	Blank Spike
ZZZZZZ	4B54438.D	08/13/15	00:49	03:43	(unrelated sample)
ZZZZZZ	4B54439.D	08/13/15	01:16	04:10	(unrelated sample)
ZZZZZZ	4B54440.D	08/13/15	01:44	04:38	(unrelated sample)
JC869-3	4B54441.D	08/13/15	02:12	05:06	(used for QC only; not part of job JC1107)
JC869-4	4B54442.D	08/13/15	02:39	05:33	(used for QC only; not part of job JC1107)
ZZZZZZ	4B54443.D	08/13/15	03:07	06:01	(unrelated sample)
ZZZZZZ	4B54444.D	08/13/15	03:34	06:28	(unrelated sample)
ZZZZZZ	4B54445.D	08/13/15	04:03	06:57	(unrelated sample)
ZZZZZZ	4B54446.D	08/13/15	04:31	07:25	(unrelated sample)
ZZZZZZ	4B54447.D	08/13/15	04:59	07:53	(unrelated sample)
ZZZZZZ	4B54448.D	08/13/15	05:27	08:21	(unrelated sample)
JC1107-8	4B54450.D	08/13/15	06:23	09:17	HSSER-RAMW06-080715
JC1107-7	4B54451.D	08/13/15	06:51	09:45	HSSER-RAMW07-080715
JC1107-9	4B54452.D	08/13/15	07:19	10:13	HSSER-RAMW05-080715
JC1107-10	4B54453.D	08/13/15	07:48	10:42	HSSER-RAMW04-080715

Instrument Performance Check (BFB)

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample:	V4B2298-BFB	Injection Date:	08/13/15
Lab File ID:	4B54454.D	Injection Time:	09:00
Instrument ID:	GCMS4B		

m/e	Ion Abundance Criteria	Raw Abundance	% Relative Abundance	Pass/Fail
50	14.99 - 40.0% of mass 95	17050	18.8	Pass
75	30.0 - 60.0% of mass 95	41786	46.1	Pass
95	Base peak, 100% relative abundance	90570	100.0	Pass
96	5.0 - 9.0% of mass 95	6316	6.97	Pass
173	Less than 2.0% of mass 174	0	0.00	(0.00) ^a Pass
174	50.0 - 120.0% of mass 95	92250	101.9	Pass
175	5.0 - 9.0% of mass 174	7549	8.33	(8.18) ^a Pass
176	95.0 - 101.0% of mass 174	90642	100.1	(98.3) ^a Pass
177	5.0 - 9.0% of mass 176	6125	6.76	(6.76) ^b Pass

(a) Value is % of mass 174

(b) Value is % of mass 176

This check applies to the following Samples, MS, MSD, Blanks, and Standards:

Lab Sample ID	Lab File ID	Date Analyzed	Time Analyzed	Hours Lapsed	Client Sample ID
V4B2298-CC2289	4B54455.D	08/13/15	09:46	00:46	Continuing cal 20
V4B2297-MB2	4B54457.D	08/13/15	10:43	01:43	Method Blank
JC869-3DUP	4B54458.D	08/13/15	11:17	02:17	Duplicate
JC869-4MS	4B54459.D	08/13/15	11:45	02:45	Matrix Spike
V4B2298-MB1	4B54461.D	08/13/15	13:09	04:09	Method Blank
V4B2298-BS	4B54462.D	08/13/15	13:38	04:38	Blank Spike
ZZZZZZ	4B54463.D	08/13/15	14:12	05:12	(unrelated sample)
JC1045-15	4B54464.D	08/13/15	14:40	05:40	(used for QC only; not part of job JC1107)
ZZZZZZ	4B54465.D	08/13/15	15:07	06:07	(unrelated sample)
ZZZZZZ	4B54466.D	08/13/15	15:36	06:36	(unrelated sample)
JC1045-15MS	4B54467.D	08/13/15	16:04	07:04	Matrix Spike
JC1045-15MSD	4B54468.D	08/13/15	16:32	07:32	Matrix Spike Duplicate
ZZZZZZ	4B54470.D	08/13/15	17:28	08:28	(unrelated sample)
ZZZZZZ	4B54471.D	08/13/15	17:56	08:56	(unrelated sample)
ZZZZZZ	4B54472.D	08/13/15	18:23	09:23	(unrelated sample)
ZZZZZZ	4B54473.D	08/13/15	18:51	09:51	(unrelated sample)
ZZZZZZ	4B54474.D	08/13/15	19:19	10:19	(unrelated sample)
ZZZZZZ	4B54475.D	08/13/15	19:48	10:48	(unrelated sample)
ZZZZZZ	4B54476.D	08/13/15	20:15	11:15	(unrelated sample)

Volatile Internal Standard Area Summary

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Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Check Std:	V4B2296-CC2289	Injection Date:	08/12/15
Lab File ID:	4B54407.D	Injection Time:	10:09
Instrument ID:	GCMS4B	Method:	SW846 8260C

	IS 1 AREA	RT	IS 2 AREA	RT	IS 3 AREA	RT	IS 4 AREA	RT	IS 5 AREA	RT
Check Std	140756	6.58	435392	8.56	495542	9.42	438556	12.60	244012	15.14
Upper Limit ^a	281512	7.08	870784	9.06	991084	9.92	877112	13.10	488024	15.64
Lower Limit ^b	70378	6.08	217696	8.06	247771	8.92	219278	12.10	122006	14.64

Lab Sample ID	IS 1 AREA	RT	IS 2 AREA	RT	IS 3 AREA	RT	IS 4 AREA	RT	IS 5 AREA	RT
V4B2296-MB	159324	6.58	450495	8.56	509837	9.42	441074	12.60	215904	15.14
V4B2296-BS	125516	6.58	422985	8.56	486014	9.42	432148	12.60	252161	15.14
ZZZZZZ	177760	6.57	440406	8.56	500506	9.42	446169	12.60	227577	15.14
ZZZZZZ	174695	6.58	446349	8.56	503883	9.42	442977	12.60	227518	15.14
ZZZZZZ	145577	6.58	442931	8.56	498816	9.42	445606	12.60	229559	15.14
JC1107-3	146811	6.58	443287	8.56	502337	9.42	441128	12.60	222699	15.14
JC1107-3MS	145924	6.58	413150	8.56	475209	9.42	424091	12.60	251934	15.14
JC1107-3MSD	152325	6.58	431587	8.56	496307	9.42	441468	12.60	260200	15.14
JC1107-11	140434	6.58	450428	8.56	498750	9.42	437102	12.60	216722	15.14
JC1107-12	137119	6.58	445018	8.56	501240	9.42	439157	12.60	212599	15.14
JC1107-1	138307	6.58	439595	8.56	488233	9.42	433020	12.60	211888	15.14
JC1107-2	140461	6.58	437134	8.56	492560	9.42	432075	12.60	213576	15.14
JC1107-4	150775	6.58	439086	8.56	491241	9.42	432149	12.60	214304	15.14
JC1107-5	149773	6.58	427592	8.56	484870	9.42	426026	12.60	211603	15.14
JC1107-6	145128	6.58	431535	8.56	484038	9.42	425202	12.60	215693	15.14

IS 1 = Tert Butyl Alcohol-D9

IS 2 = Pentafluorobenzene

IS 3 = 1,4-Difluorobenzene

IS 4 = Chlorobenzene-D5

IS 5 = 1,4-Dichlorobenzene-d4

(a) Upper Limit = + 100% of check standard area; Retention time + 0.5 minutes.

(b) Lower Limit = -50% of check standard area; Retention time -0.5 minutes.

Volatile Internal Standard Area Summary

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Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Check Std:	V4B2297-CC2289	Injection Date:	08/12/15
Lab File ID:	4B54431.D	Injection Time:	21:35
Instrument ID:	GCMS4B	Method:	SW846 8260C

	IS 1 AREA	RT	IS 2 AREA	RT	IS 3 AREA	RT	IS 4 AREA	RT	IS 5 AREA	RT
Check Std	129429	6.58	415468	8.56	477620	9.42	432693	12.60	251516	15.14
Upper Limit ^a	258858	7.08	830936	9.06	955240	9.92	865386	13.10	503032	15.64
Lower Limit ^b	64715	6.08	207734	8.06	238810	8.92	216347	12.10	125758	14.64

Lab Sample ID	IS 1 AREA	RT	IS 2 AREA	RT	IS 3 AREA	RT	IS 4 AREA	RT	IS 5 AREA	RT
V4B2297-MB	149010	6.58	437268	8.56	487916	9.42	429546	12.60	212618	15.14
V4B2297-BS	119545	6.57	416714	8.56	479398	9.42	432369	12.60	251910	15.14
ZZZZZZ	149580	6.58	445004	8.56	502390	9.42	436916	12.60	215385	15.14
ZZZZZZ	141252	6.58	435800	8.56	486562	9.42	427224	12.60	209932	15.14
ZZZZZZ	136795	6.57	434712	8.56	483983	9.42	430219	12.60	208221	15.14
JC869-3	144272	6.58	433663	8.56	489597	9.42	429586	12.60	208261	15.14
JC869-4	137817	6.58	429129	8.56	483587	9.42	426165	12.60	207686	15.14
ZZZZZZ	140668	6.58	429440	8.56	488300	9.42	427955	12.60	208028	15.14
ZZZZZZ	142053	6.57	429641	8.56	482462	9.42	422562	12.60	204634	15.14
ZZZZZZ	147189	6.58	428108	8.56	479980	9.42	421440	12.60	202976	15.14
ZZZZZZ	143075	6.58	423446	8.56	479173	9.42	429412	12.60	222190	15.14
ZZZZZZ	129439	6.58	429498	8.56	485670	9.42	425477	12.60	202610	15.14
ZZZZZZ	137442	6.57	428832	8.56	483312	9.42	440691	12.60	252677	15.14
JC1107-8	144786	6.58	464734	8.56	523670	9.42	454873	12.60	216894	15.14
JC1107-7	181708	6.58	443640	8.56	499255	9.42	435289	12.60	212248	15.14
JC1107-9	142504	6.58	438675	8.56	494969	9.42	432890	12.60	208348	15.14
JC1107-10	134311	6.58	433187	8.56	481637	9.42	421297	12.60	205804	15.14

IS 1 = Tert Butyl Alcohol-D9

IS 2 = Pentafluorobenzene

IS 3 = 1,4-Difluorobenzene

IS 4 = Chlorobenzene-D5

IS 5 = 1,4-Dichlorobenzene-d4

(a) Upper Limit = + 100% of check standard area; Retention time + 0.5 minutes.

(b) Lower Limit = -50% of check standard area; Retention time -0.5 minutes.

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Volatile Internal Standard Area Summary

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Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Check Std:	V4B2298-CC2289	Injection Date:	08/13/15
Lab File ID:	4B54455.D	Injection Time:	09:46
Instrument ID:	GCMS4B	Method:	SW846 8260C

	IS 1 AREA	RT	IS 2 AREA	RT	IS 3 AREA	RT	IS 4 AREA	RT	IS 5 AREA	RT
Check Std	150330	6.57	418083	8.56	472304	9.41	415435	12.60	226426	15.14
Upper Limit ^a	300660	7.07	836166	9.06	944608	9.91	830870	13.10	452852	15.64
Lower Limit ^b	75165	6.07	209042	8.06	236152	8.91	207718	12.10	113213	14.64

Lab Sample ID	IS 1 AREA	RT	IS 2 AREA	RT	IS 3 AREA	RT	IS 4 AREA	RT	IS 5 AREA	RT
V4B2297-MB2	129837	6.58	435467	8.56	485700	9.42	421308	12.60	200871	15.14
JC869-3DUP	129185	6.58	418519	8.56	473275	9.42	402037	12.60	185731	15.14
JC869-4MS	137687	6.58	412721	8.56	476155	9.42	427821	12.60	246362	15.14
V4B2298-MB1	141611	6.58	423356	8.56	470956	9.41	406520	12.60	189367	15.14
V4B2298-BS	133396	6.57	415006	8.56	473437	9.42	421683	12.60	239628	15.14
ZZZZZZ	153534	6.58	428237	8.56	479142	9.42	408818	12.60	199520	15.14
JC1045-15	139403	6.58	433083	8.56	486646	9.42	425186	12.60	205509	15.14
ZZZZZZ	145344	6.58	432936	8.56	485272	9.42	420337	12.60	199487	15.14
ZZZZZZ	141731	6.58	425207	8.56	470874	9.42	412527	12.60	204040	15.14
JC1045-15MS	134272	6.58	405035	8.56	462082	9.42	418249	12.60	241503	15.14
JC1045-15MSD	136201	6.58	418290	8.56	479337	9.42	430460	12.60	248080	15.14
ZZZZZZ	141097	6.58	437011	8.56	487946	9.42	428097	12.60	205837	15.14
ZZZZZZ	143967	6.58	431844	8.56	485058	9.42	424528	12.60	212720	15.14
ZZZZZZ	144271	6.58	429633	8.56	483460	9.42	423107	12.60	201687	15.14
ZZZZZZ	141731	6.58	425547	8.56	478561	9.42	416690	12.60	199888	15.14
ZZZZZZ	144549	6.58	427563	8.56	477656	9.42	416181	12.60	200991	15.14
ZZZZZZ	141741	6.58	419551	8.56	471047	9.42	417921	12.60	198968	15.14
ZZZZZZ	148549	6.58	425269	8.56	479509	9.42	418275	12.60	201366	15.14

IS 1 = Tert Butyl Alcohol-D9

IS 2 = Pentafluorobenzene

IS 3 = 1,4-Difluorobenzene

IS 4 = Chlorobenzene-D5

IS 5 = 1,4-Dichlorobenzene-d4

(a) Upper Limit = + 100% of check standard area; Retention time + 0.5 minutes.

(b) Lower Limit = -50% of check standard area; Retention time -0.5 minutes.

Volatile Surrogate Recovery Summary

Page 1 of 1

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Method: SW846 8260C

Matrix: AQ

Samples and QC shown here apply to the above method

Lab Sample ID	Lab File ID	S1	S2	S3	S4
JC1107-1	4B54425.D	97	92	100	105
JC1107-2	4B54426.D	97	94	99	104
JC1107-3	4B54416.D	98	92	99	104
JC1107-4	4B54427.D	97	93	99	104
JC1107-5	4B54428.D	98	94	99	104
JC1107-6	4B54429.D	98	94	99	103
JC1107-7	4B54451.D	97	92	98	104
JC1107-8	4B54450.D	97	91	98	107
JC1107-9	4B54452.D	97	93	98	106
JC1107-10	4B54453.D	98	94	100	105
JC1107-11	4B54423.D	96	91	100	104
JC1107-12	4B54424.D	96	92	99	105
JC1107-3MS	4B54420.D	98	94	98	95
JC1107-3MSD	4B54421.D	99	94	99	95
JC869-3DUP	4B54458.D	99	92	97	109
JC869-4MS	4B54459.D	101	95	99	97
V4B2296-BS	4B54410.D	100	94	100	96
V4B2296-MB	4B54409.D	98	92	98	105
V4B2297-BS	4B54434.D	100	95	100	96
V4B2297-MB	4B54433.D	98	93	98	104
V4B2297-MB2	4B54457.D	96	92	99	106

Surrogate Compounds	Recovery Limits
---------------------	-----------------

S1 = Dibromofluoromethane	76-120%
S2 = 1,2-Dichloroethane-D4	73-122%
S3 = Toluene-D8	84-119%
S4 = 4-Bromofluorobenzene	78-117%

Raw Data: 4B54242.D 4B54243.D 4B54244.D 4B54245.D 4B54246.D 4B54247.D 4B54248.D 4B54249.D
 4B54250.D 4B54251.D

Initial Calibration Summary

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Page 1 of 5

Sample: V4B2289-ICC2289
 Lab FileID: 4B54249.D

Response Factor Report MS4B

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)

Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um

Last Update : Mon Aug 10 08:19:32 2015

Response via : Initial Calibration

Calibration Files

1	=4B54244.D	0.5	=4B54243.D	100	=4B54250.D	50	=4B54249.D
20	=4B54248.D	200	=4B54251.D	5	=4B54246.D	2	=4B54245.D
10	=4B54247.D	0.2	=4B54242.D		=		=

Compound

	1	0.5	100	50	20	200	5	2	10	0.2	Avg	%RSD
<hr/>												
1)	tert butyl alcohol-d9											
2)	tertiary butyl alcohol	1.351		1.204	1.128	1.249	1.143	1.152	1.209	1.241	1.209	6.02
3)	1,4-dioxane			0.122	0.113	0.120	0.119	0.106	0.092	0.111	0.112	9.46
4)	Ethanol										0.000#	-1.00
5)	I pentafluorobenzene											
6)	CHLOROTRIFLUOROETHENE										0.000#	-1.00
7)	chlorodifluoromethane	0.575		0.504	0.479	0.486	0.505	0.472	0.527	0.496	0.506	6.48
8)	dichlorodifluoromethane			0.504	0.474	0.460	0.477	0.497	0.474	0.504	0.484	3.59
9)	Freon 114										0.000#	-1.00
10)	chloromethane	0.181	0.188	0.193	0.176	0.174	0.193	0.188	0.191	0.182	0.185	3.86
11)	vinyl chloride	0.497	0.466	0.571	0.527	0.518	0.552	0.545	0.562	0.551	0.532	6.36
12)	bromomethane	0.285	0.344	0.307	0.300	0.289	0.253	0.315	0.338	0.317	0.305	9.14
13)	chloroethane	0.189	0.211	0.221	0.210	0.207	0.193	0.220	0.214	0.219	0.209	5.55
14)	vinyl bromide										0.000#	-1.00
15)	trichlorofluoromethane	0.391		0.548	0.508	0.491	0.514	0.524	0.514	0.531	0.503	9.55
16)	1,3-butadiene										0.000#	-1.00
17)	Pentane										0.000#	-1.00
18)	freon 123a										0.000#	-1.00
19)	ethyl ether	0.208	0.158	0.198	0.188	0.188	0.201	0.183	0.198	0.193	0.191	7.49
20)	2-chloropropane--The compound does not meet initial criteria.											
		0.055	0.056	0.066	0.047	0.066		0.073			0.061	15.71
21)	acrolein	0.061	0.058	0.071	0.067	0.062	0.070	0.061	0.059	0.065	0.064	7.35
22)	1,1-dichloroethene	0.457	0.323	0.386	0.366	0.380	0.404	0.392	0.430	0.415	0.395	9.75
23)	acetone											

Initial Calibration Summary

Page 2 of 5

Job Number: JC1107

Sample: V4B2289-ICC2289
Lab FileID: 4B54249.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

	0.035 0.034 0.033 0.034 0.033 0.033 0.031	0.033	3.62
24)	allyl chloride 0.320 0.311 0.270 0.258 0.322 0.255 0.260 0.318	0.289	10.69
25)	acetonitrile 0.032 0.031 0.034 0.031 0.029 0.031 0.035	0.032	5.56
26)	iodomethane 0.778 0.601 0.768 0.734 0.729 0.785 0.740 0.775 0.800	0.746	7.98
27)	carbon disulfide 1.249 0.924 1.171 1.107 1.158 1.187 1.143 1.246 1.243	1.159	8.72
28)	methylene chloride 0.434 0.358 0.425 0.413 0.407 0.434 0.413 0.431 0.447	0.418	6.16
29)	methyl acetate 0.049 0.062 0.060 0.059 0.064 0.048 0.051 0.058	0.056	10.68
30)	1-chloropropane 0.680 0.652 0.668 0.669 0.751 0.822 0.752	0.714	8.79
31)	methyl tert butyl ether 1.130 0.893 1.075 1.034 1.043 1.069 1.057 1.084 1.097	1.053	6.34
32)	trans-1,2-dichloroethene 0.478 0.371 0.368 0.345 0.359 0.376 0.383 0.408 0.407	0.388	10.19
33)	di-isopropyl ether 1.503 1.187 1.394 1.367 1.373 1.353 1.353 1.484 1.420	1.381	6.61
34)	2-butanone 0.048 0.043 0.044 0.047 0.040	0.043	6.65
35)	1,1-dichloroethane 0.739 0.713 0.683 0.718 0.702 0.734 0.773 0.776	0.730	4.47
36)	chloroprene 0.605 0.531 0.499 0.525 0.520 0.536 0.556 0.551	0.540	5.87
37)	acrylonitrile 0.160 0.152 0.148 0.146 0.153 0.145 0.135 0.151	0.149	4.94
38)	vinyl acetate 0.050 0.058 0.056 0.057 0.059 0.051 0.066 0.054	0.056	8.84
39)	ethyl tert-butyl ether 1.281 1.073 1.258 1.227 1.216 1.249 1.179 1.318 1.243	1.227	5.70
40)	ethyl acetate 0.061 0.059 0.058 0.059 0.057 0.065 0.062	0.060	4.48
41)	2,2-dichloropropane 0.359 0.272 0.272 0.302 0.244 0.320 0.362 0.338	0.309	14.11
42)	cis-1,2-dichloroethene 0.455 0.419 0.399 0.399 0.421 0.401 0.447 0.433	0.422	5.21
43)	methylacrylate 0.061 0.057 0.056 0.062 0.051	0.056	7.24
44)	propionitrile 0.061 0.051 0.063 0.059 0.060 0.063 0.061 0.058 0.063	0.060	6.51
45)	bromochloromethane 0.212 0.224 0.214 0.213 0.229 0.206 0.212 0.223	0.216	3.51
46)	tetrahydrofuran 0.161 0.141 0.134 0.142 0.136 0.143 0.151 0.145	0.144	6.08
47)	chloroform 0.499 0.364 0.431 0.425 0.425 0.440 0.435 0.476 0.470 0.402	0.437	8.79
48)	T-BUTYL FORMATE 0.261 0.307 0.297 0.279 0.311 0.251 0.259 0.268	0.279	8.33
49)	dibromofluoromethane (s) 0.358 0.353 0.368 0.363 0.368 0.368 0.360 0.348 0.366 0.363	0.362	1.90
50)	1,2-dichloroethane-d4 (s) 0.398 0.399 0.391 0.397 0.404 0.380 0.402 0.401 0.399 0.403	0.397	1.83
51)	freon 113 0.242 0.232 0.227 0.236 0.233 0.250 0.241 0.256	0.240	4.08
52)	methacrylonitrile 0.273 0.243 0.273 0.256 0.259 0.269 0.261 0.276 0.272	0.265	4.01
53)	1,1,1-trichloroethane		

Initial Calibration Summary

Page 3 of 5

Job Number: JC1107

Sample: V4B2289-ICC2289
Lab FileID: 4B54249.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

54)	cyclohexane	0.472 0.534	0.457 0.435 0.455 0.448 0.449 0.480 0.498 0.500 0.460 0.481 0.515 0.501 0.517 0.539	0.462 0.506	4.41 5.21
55)	iso-butyl alcohol		0.017 0.015 0.016 0.017 0.018	0.017	6.13
56)	I 1,4-difluorobenzene		-----ISTD-----		
57)	epichlorohydrin	0.039 0.039	0.028 0.036 0.034 0.031 0.036 0.036 0.032 0.036	0.034	9.51
58)	n-butyl alcohol	0.011 0.011	0.009 0.010 0.009 0.010 0.010 0.010 0.009 0.010	0.010#	6.67
59)	carbon tetrachloride	0.422 0.422	0.395 0.372 0.389 0.396 0.389 0.413 0.424	0.400	4.52
60)	1,1-dichloropropene	0.479 0.479	0.430 0.403 0.433 0.429 0.444 0.466 0.473	0.445	5.88
61)	hexane	0.389 0.389	0.320 0.298 0.317 0.313 0.354 0.358 0.352	0.338	9.00
62)	Tert Amyl alcohol			0.000#	-1.00
63)	benzene	1.477 1.477	1.164 1.263 1.211 1.251 1.265 1.310 1.422 1.378 1.345	1.309	7.44
64)	iso-octane	1.009 1.009	0.959 1.013 0.995 1.131 1.143 1.107	1.051	7.03
65)	tert-amyl methyl ether	0.203 0.203	0.197 0.205 0.205 0.204 0.204 0.235 0.213	0.208	5.57
66)	heptane	0.249 0.249	0.225 0.210 0.225 0.221 0.246 0.258 0.226	0.232	7.10
67)	isopropyl acetate	0.140 0.140	0.139 0.147 0.139 0.202	0.173	16.45
68)	1,2-dichloroethane	0.477 0.477	0.339 0.420 0.423 0.435 0.404 0.441 0.471 0.472	0.431	9.99
69)	trichloroethene	0.384 0.384	0.347 0.332 0.334 0.348 0.348 0.354 0.371	0.352	4.99
70)	Tert-amyl Ethyl Ether			0.000#	-1.00
71)	ethyl acrylate	0.495 0.495	0.478 0.457 0.478 0.625 0.508 0.501	0.506	10.93
72)	2-nitropropane	0.160 0.160	0.119 0.139 0.136 0.140 0.131 0.145 0.151 0.150	0.141	8.59
73)	2-chloroethyl vinyl ether	0.219 0.219	0.187 0.218 0.215 0.211 0.213 0.207 0.218 0.218 0.180	0.209	6.65
74)	methyl methacrylate	0.082 0.082	0.089 0.085 0.084 0.091 0.083 0.078 0.082	0.084	4.85
75)	1,2-dichloropropane	0.381 0.381	0.367 0.361 0.370 0.360 0.365 0.404 0.401	0.376	4.66
76)	dibromomethane	0.239 0.239	0.175 0.226 0.220 0.217 0.227 0.218 0.213 0.235	0.219	8.53
77)	methylcyclohexane	0.464 0.464	0.456 0.416 0.439 0.450 0.470 0.461 0.467	0.453	3.97
78)	bromodichloromethane	0.496 0.496	0.356 0.486 0.472 0.465 0.487 0.480 0.495 0.498	0.471	9.40
79)	cis-1,3-dichloropropene	0.623 0.623	0.615 0.590 0.591 0.612 0.597 0.620 0.631	0.610	2.53
80)	toluene-d8 (s)	1.178 1.178	1.175 1.178 1.185 1.177 1.168 1.173 1.179 1.186 1.174	1.177	0.44
81)	4-methyl-2-pentanone	0.139 0.139	0.148 0.141 0.147 0.148 0.151 0.134 0.148	0.144	4.03
82)	toluene	0.908 0.908	0.816 0.771 0.784 0.833 0.828 0.873 0.860	0.834	5.45
83)	3-methyl-1-butanol				

Initial Calibration Summary

Page 4 of 5

Job Number: JC1107

Sample: V4B2289-ICC2289
Lab FileID: 4B54249.DAccount: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

0.018	0.016	0.015	0.016	0.015	0.017	0.017	0.016	0.016	6.45
84) trans-1,3-dichloropropene	0.554	0.415	0.557	0.542	0.537	0.551	0.551	0.535	0.564
85) ethyl methacrylate	0.430	0.477	0.449	0.428	0.477	0.421	0.409	0.445	0.442
86) 1,1,2-trichloroethane	0.306	0.294	0.283	0.274	0.293	0.278	0.300	0.294	0.290
87) 2-hexanone	0.135	0.142	0.127	0.128	0.138	0.138	0.117	0.131	0.132
88) I chlorobenzene-d5									ISTD-----
89) tetrachloroethene	0.458	0.417	0.393	0.418	0.414	0.449	0.469	0.457	0.434
90) 1,3-dichloropropane	0.649	0.453	0.587	0.576	0.581	0.583	0.619	0.633	0.616
91) butyl acetate	0.275	0.271	0.258	0.259	0.274	0.276	0.258	0.277	0.268
92) 3,3-DIMETHYL-1-BUTANOL	0.043	0.035	0.038	0.036	0.040	0.038	0.042	0.042	0.041
93) dibromochloromethane	0.480	0.468	0.444	0.439	0.479	0.443	0.464	0.464	0.460
94) 1,2-dibromoethane	0.411	0.309	0.404	0.382	0.379	0.413	0.389	0.389	0.394
95) n-butyl ether									0.000# -1.00
96) chlorobenzene	1.177	0.836	1.060	1.001	1.028	1.086	1.057	1.168	1.105
97) 1,1,1,2-tetrachloroethane	0.424	0.396	0.385	0.391	0.408	0.403	0.405	0.421	0.404
98) ethylbenzene	1.982	1.384	1.751	1.665	1.721	1.773	1.801	1.919	1.856
99) m,p-xylene	0.759	0.520	0.682	0.639	0.654	0.699	0.675	0.707	0.706
100) o-xylene	0.730	0.504	0.709	0.665	0.667	0.724	0.686	0.694	0.710
101) styrene	1.165	1.199	1.124	1.126	1.226	1.102	1.116	1.186	1.155
102) bromoform	0.367	0.377	0.355	0.344	0.388	0.344	0.332	0.355	0.358
103) I 1,4-dichlorobenzene-d									ISTD-----
104) isopropylbenzene	3.596	3.029	2.849	2.979	3.113	3.176	3.336	3.314	3.174
105) 4-bromofluorobenzene (s)	0.877	0.889	0.852	0.854	0.867	0.857	0.880	0.881	0.867
106) cyclohexanone--The compound does not meet initial criteria.	0.092	0.078	0.092	0.073	0.106	0.094	0.099	0.091	12.79
107) bromobenzene	1.044	0.764	0.926	0.875	0.897	0.963	0.953	0.985	0.973
108) 1,1,2,2-tetrachloroethane	0.982	0.794	0.926	0.891	0.906	0.944	0.966	0.981	0.975
109) trans-1,4-dichloro-2-butene	0.261	0.194	0.248	0.232	0.235	0.251	0.251	0.223	0.257
110) 1,2,3-trichloropropane	0.214	0.179	0.220	0.214	0.228	0.221	0.239	0.228	0.239
111) n-propylbenzene	4.087	3.067	3.493	3.255	3.426	3.565	3.703	3.872	3.822
112) 4-ETHYLTOLEUENE									0.000# -1.00
113) 2-chlorotoluene									

Initial Calibration Summary

Page 5 of 5

Job Number: JC1107

Sample: V4B2289-ICC2289
Lab FileID: 4B54249.DAccount: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

114)	4-chlorotoluene	0.879 0.789 0.737 0.769 0.816 0.798 0.864 0.854	0.813	6.08
115)	1,3,5-trimethylbenzene	2.629 1.912 2.281 2.163 2.218 2.348 2.358 2.520 2.450	2.320	9.12
116)	tert-butylbenzene	2.946 2.581 2.431 2.523 2.617 2.644 2.848 2.785	2.672	6.50
117)	pentachloroethane	2.577 2.142 2.261 2.144 2.210 2.291 2.334 2.467 2.434	2.318	6.47
118)	1,2,4-trimethylbenzene	0.524 0.406 0.462 0.448 0.455 0.507 0.488 0.537 0.490	0.480	8.58
119)	sec-butylbenzene	2.803 2.082 2.536 2.442 2.511 2.592 2.636 2.770 2.792	2.574	8.77
120)	1,3-dichlorobenzene	3.819 3.417 3.224 3.342 3.490 3.502 3.692 3.689	3.522	5.67
121)	p-isopropyltoluene	1.908 1.417 1.706 1.639 1.659 1.757 1.731 1.833 1.830 1.468	1.695	9.25
122)	1,4-dichlorobenzene	3.116 2.887 2.731 2.830 2.915 2.943 3.055 3.075	2.944	4.47
123)	benzyl chloride	1.901 1.471 1.694 1.620 1.635 1.746 1.689 1.841 1.777 1.578	1.695	7.52
124)	1,2-dichlorobenzene	1.897 1.673 1.672 1.633 1.603 1.686 1.702 1.740 1.644	1.694	5.06
125)	1,4-DIETHYLBENZENE	1.862 1.320 1.666 1.606 1.632 1.681 1.658 1.712 1.757	1.655	8.88
			0.000#	-1.00
126)	n-butylbenzene	1.500 1.482 1.387 1.428 1.484 1.416 1.441 1.544	1.460	3.52
127)	1,2,4,5-TETRAMETHYLBENZENE		0.000#	-1.00
128)	1,2-dibromo-3-chloropropane	0.162 0.150 0.140 0.145 0.152 0.144 0.146 0.155	0.149	4.78
129)	1,3,5-TRICHLOROBENZENE	1.375 1.058 1.343 1.305 1.318 1.352 1.323 1.401 1.404	1.320	7.91
130)	1,2,4-trichlorobenzene	1.032 1.117 1.068 1.037 1.154 0.990 0.997 1.063	1.057	5.34
131)	hexachlorobutadiene	0.575 0.581 0.640 0.595 0.674 0.713 0.725	0.643	9.70
132)	naphthalene	1.641 2.051 1.913 1.867 2.165 1.712 1.677 1.841	1.858	9.88
133)	1,2,3-trichlorobenzene	0.838 0.933 0.895 0.908 0.971 0.841 0.846 0.902	0.892	5.36
134)	hexachloroethane	0.653 0.598 0.562 0.580 0.623 0.606 0.622 0.640	0.610	4.95

(##) = Out of Range ### Number of calibration levels exceeded format ###

M4B2289.M Mon Aug 10 08:20:19 2015 GCMS4B

Initial Calibration Verification

Job Number: JC1107

Sample: V4B2289-ICV2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54254.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Evaluate Continuing Calibration Report

Data File : C:\MSDCHEM\1\DATA\4B54254.D Vial: 14
 Acq On : 5 Aug 2015 3:20 pm Operator: TOANP
 Sample : icv2289-50 Inst : MS4B
 Misc : MS88663,V4B2289,w,,,1 Multiplr: 1.00
 MS Integration Params: RTEINT.P

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 Last Update : Mon Aug 10 08:19:32 2015
 Response via : Multiple Level Calibration

Min. RRF : 0.010 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 200%

	Compound	TrueValue	AvgRF	CCRF	%Dev	Area*	Dev(min)	R.T.
1	tert butyl alcohol-d9	500.00	1.000	1.000	0.0	100	0.00	6.57
2 M	tertiary butyl alcohol	250.00	1.209	1.239	-2.5	110	0.00	6.67
3 M	1,4-dioxane	1250.00	0.112	0.126	-12.5	111	0.00	10.11
4	Ethanol		-----NA-----					
5 I	pentafluorobenzene	50.00	1.000	1.000	0.0	103	0.00	8.56
6	CHLOROTRIFLUOROETHENE		-----NA-----					
7 M	chlorodifluoromethane	50.00	0.506	0.448	11.5	97	0.00	3.63
8 M	dichlorodifluoromethane	50.00	0.484	0.379	21.7	83	-0.01	3.60
9	Freon 114		-----NA-----					
10 M	chloromethane	50.00	0.185	0.170	8.1	100	0.00	3.96
11 M	vinyl chloride	50.00	0.532	0.467	12.2	91	0.00	4.17
12 M	bromomethane	50.00	0.305	0.309	-1.3	106	0.01	4.76
13 M	chloroethane	50.00	0.209	0.243	-16.3	119	0.01	4.92
14	vinyl bromide		-----NA-----					
15 M	trichlorofluoromethane	50.00	0.503	0.463	8.0	94	0.00	5.27
16	1,3-butadiene		-----NA-----					
17	Pentane		-----NA-----					
18	freon 123a		-----NA-----					
19 M	ethyl ether	50.00	0.191	0.190	0.5	105	0.00	5.63
20	2-chloropropane	50.00	0.061	0.040	34.4#	74	0.00	5.81
21 M	acrolein	500.00	0.064	0.080	-25.0	123	0.00	5.86
22 M	1,1-dichloroethene	50.00	0.395	0.387	2.0	109	0.01	6.03
23 M	acetone	50.00	0.033	0.036	-9.1	109	0.00	6.05
24 M	allyl chloride	50.00	0.289	0.369	-27.7	141	-0.01	6.45
25 M	acetonitrile	500.00	0.032	0.031	3.1	103	0.00	6.45
26 M	iodomethane	50.00	0.746	0.687	7.9	97	0.01	6.28
27 M	carbon disulfide	50.00	1.159	1.081	6.7	101	0.00	6.39
28 M	methylene chloride	50.00	0.418	0.421	-0.7	105	0.01	6.65
29 M	methyl acetate	50.00	0.056	0.055	1.8	96	0.00	6.44
30	1-chloropropane	50.00	0.714	0.650	9.0	103	0.00	6.65
31 M	methyl tert butyl ether	100.00	1.053	0.997	5.3	100	0.00	6.89
32 M	trans-1,2-dichloroethene	50.00	0.388	0.347	10.6	104	0.00	6.95
33 M	di-isopropyl ether	50.00	1.381	1.334	3.4	101	0.00	7.41
34 M	2-butanone	50.00	0.044	0.045	-2.3	106	0.00	8.08
35 M	1,1-dichloroethane	50.00	0.730	0.690	5.5	104	0.00	7.45
36 M	chloroprene	50.00	0.540	0.487	9.8	101	0.00	7.54
37 M	acrylonitrile	250.00	0.149	0.151	-1.3	106	0.00	6.92
38 M	vinyl acetate	50.00	0.056	0.061	-8.9	113	0.00	7.43
39 M	ethyl tert-butyl ether	50.00	1.227	1.211	1.3	102	0.00	7.83
40 M	ethyl acetate	50.00	0.060	0.057	5.0	99	0.00	8.08
41 M	2,2-dichloropropane	50.00	0.309	0.240	22.3	91	0.00	8.10

6.92

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Initial Calibration Verification

Page 2 of 3

Job Number: JC1107

Sample: V4B2289-ICV2289
Lab FileID: 4B54254.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

42 M	cis-1,2-dichloroethene	50.00	0.422	0.397	5.9	103	0.00	8.11
43	methylacrylate	50.00	0.057	0.057	0.0	103	0.00	8.16
44 M	propionitrile	500.00	0.060	0.057	5.0	99	0.00	8.18
45 M	bromochloromethane	50.00	0.216	0.214	0.9	103	0.00	8.40
46 M	tetrahydrofuran	50.00	0.144	0.136	5.6	104	0.00	8.43
47 M	chloroform	50.00	0.437	0.427	2.3	104	0.00	8.44
48 M	T-BUTYL FORMATE	50.00	0.279	0.280	-0.4	98	0.00	8.46
49 S	dibromofluoromethane (s)	50.00	0.362	0.362	0.0	103	0.00	8.63
50 S	1,2-dichloroethane-d4 (s)	50.00	0.397	0.379	4.5	99	0.00	9.02
51 M	freon 113	50.00	0.240	0.222	7.5	101	0.00	5.97
52 M	methacrylonitrile	50.00	0.265	0.259	2.3	105	0.00	8.34
53 M	1,1,1-trichloroethane	50.00	0.462	0.418	9.5	99	0.00	8.67
54 M	cyclohexane	50.00	0.506	0.393	22.3	88	0.00	8.73
55	iso-butyl alcohol	500.00	0.016	0.016	0.0	107	0.00	8.82
56 I	1,4-difluorobenzene	50.00	1.000	1.000	0.0	103	0.00	9.42
57 M	epichlorohydrin	250.00	0.034	0.035	-2.9	105	0.00	10.63
58 M	n-butyl alcohol	2500.00	0.010	0.009#	10.0	108	0.00	9.53
59 M	carbon tetrachloride	50.00	0.400	0.353	11.8	98	0.00	8.86
60 M	1,1-dichloropropene	50.00	0.445	0.420	5.6	107	0.00	8.83
61 M	hexane	50.00	0.338	0.365	-8.0	127	0.00	7.20
62	Tert Amyl alcohol				-----NA-----			
63 M	benzene	50.00	1.309	1.226	6.3	104	0.00	9.08
64 m	iso-octane	50.00	1.051	0.739	29.7	79	0.00	9.07
65 M	tert-amyl methyl ether	50.00	0.208	0.193	7.2	97	0.00	9.09
66 M	heptane	50.00	0.232	0.243	-4.7	119	0.00	9.21
67 M	isopropyl acetate	50.00	0.156	0.135	13.5	100	0.00	8.99
68 M	1,2-dichloroethane	50.00	0.431	0.409	5.1	100	0.00	9.10
69 M	trichloroethene	50.00	0.352	0.332	5.7	103	0.00	9.74
70	Tert-amyl Ethyl Ether				-----NA-----			
71	ethyl acrylate	50.00	0.506	0.464	8.3	100	0.00	9.73
72 M	2-nitropropane	50.00	0.141	0.139	1.4	105	0.00	10.50
73 M	2-chloroethyl vinyl ether	250.00	0.209	0.233	-11.5	112	0.00	10.50
74 M	methyl methacrylate	50.00	0.084	0.089	-6.0	108	0.00	9.99
75 M	1,2-dichloropropane	50.00	0.376	0.361	4.0	103	0.00	10.00
76 M	dibromomethane	50.00	0.219	0.213	2.7	100	0.00	10.16
77 M	methylcyclohexane	50.00	0.453	0.396	12.6	98	0.00	9.94
78 M	bromodichloromethane	50.00	0.471	0.468	0.6	102	0.00	10.28
79 M	cis-1,3-dichloropropene	50.00	0.610	0.584	4.3	102	0.00	10.72
80 S	toluene-d8 (s)	50.00	1.177	1.169	0.7	102	0.00	11.01
81 M	4-methyl-2-pentanone	50.00	0.144	0.147	-2.1	107	0.00	10.82
82 M	toluene	50.00	0.834	0.778	6.7	104	0.00	11.09
83 M	3-methyl-1-butanol	1000.00	0.016	0.015	6.3	108	0.00	10.83
84 M	trans-1,3-dichloropropene	50.00	0.534	0.501	6.2	95	0.00	11.29
85 M	ethyl methacrylate	50.00	0.442	0.443	-0.2	102	0.00	11.27
86 M	1,1,2-trichloroethane	50.00	0.290	0.282	2.8	103	0.00	11.52
87 M	2-hexanone	50.00	0.132	0.131	0.8	106	0.00	11.69
88 I	chlorobenzene-d5	50.00	1.000	1.000	0.0	101	0.00	12.60
89 M	tetrachloroethene	50.00	0.434	0.399	8.1	103	0.00	11.69
90 M	1,3-dichloropropane	50.00	0.588	0.571	2.9	100	0.00	11.71
91 M	butyl acetate	50.00	0.268	0.260	3.0	102	0.00	11.76
92 M	3,3-DIMETHYL-1-BUTANOL	500.00	0.039	0.035	10.3	99	0.00	11.87
93 M	dibromochloromethane	50.00	0.460	0.449	2.4	102	0.00	11.99
94 M	1,2-dibromoethane	50.00	0.385	0.385	0.0	102	0.00	12.15
95	n-butyl ether				-----NA-----			
96 M	chlorobenzene	50.00	1.057	1.032	2.4	104	0.00	12.63
97 M	1,1,1,2-tetrachloroethane	50.00	0.404	0.378	6.4	99	0.00	12.70
98 M	ethylbenzene	50.00	1.761	1.663	5.6	101	0.00	12.69
99 M	m,p-xylene	100.00	0.671	0.647	3.6	102	0.00	12.80

6.9.2
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Initial Calibration Verification

Page 3 of 3

Job Number: JC1107

Sample: V4B2289-ICV2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54254.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

100 M	<i>o</i> -xylene	50.00	0.677	0.677	0.0	103	0.00	13.25
101 M	styrene	50.00	1.155	1.130	2.2	102	0.00	13.27
102 M	bromoform	50.00	0.358	0.359	-0.3	102	0.00	13.57
103 I	1,4-dichlorobenzene-d4	50.00	1.000	1.000	0.0	99	0.00	15.14
104 M	isopropylbenzene	50.00	3.174	2.910	8.3	101	0.00	13.63
105 S	4-bromofluorobenzene (s)	50.00	0.873	0.863	1.1	100	0.00	13.86
106	cyclohexanone	500.00	0.091	0.063	30.8#	80	0.00	13.83
107 M	bromobenzene	50.00	0.931	0.888	4.6	100	0.00	14.07
108 M	1,1,2,2-tetrachloroethane	50.00	0.930	0.890	4.3	98	0.00	13.98
109 M	trans-1,4-dichloro-2-butene	50.00	0.239	0.228	4.6	97	0.00	14.02
110 M	1,2,3-trichloropropane	50.00	0.220	0.214	2.7	99	0.00	14.06
111 M	n-propylbenzene	50.00	3.588	3.502	2.4	106	0.00	14.08
112	4-ETHYLTOULUENE		-----NA-----					
113 M	2-chlorotoluene	50.00	0.813	0.746	8.2	100	0.00	14.24
114 M	4-chlorotoluene	50.00	2.320	2.169	6.5	99	0.00	14.35
115 M	1,3,5-trimethylbenzene	50.00	2.672	2.467	7.7	100	0.00	14.24
116 M	tert-butylbenzene	50.00	2.318	2.204	4.9	101	0.00	14.62
117 M	pentachloroethane	50.00	0.480	0.472	1.7	104	0.00	14.72
118 M	1,2,4-trimethylbenzene	50.00	2.574	2.571	0.1	104	0.00	14.68
119 M	sec-butylbenzene	50.00	3.522	3.232	8.2	99	0.00	14.87
120 M	1,3-dichlorobenzene	50.00	1.695	1.651	2.6	99	0.00	15.08
121 M	p-isopropyltoluene	50.00	2.944	2.772	5.8	100	0.00	15.00
122 M	1,4-dichlorobenzene	50.00	1.695	1.633	3.7	99	0.00	15.17
123	benzyl chloride	50.00	1.694	1.593	6.0	96	0.00	15.30
124 M	1,2-dichlorobenzene	50.00	1.655	1.636	1.1	100	0.00	15.61
125	1,4-DIETHYLBENZENE		-----NA-----					
126 M	n-butylbenzene	50.00	1.460	1.397	4.3	99	0.00	15.46
127	1,2,4,5-TETRAMETHYLBENZEN		-----NA-----					
128 M	1,2-dibromo-3-chloropropane	50.00	0.149	0.147	1.3	104	0.00	16.46
129	1,3,5-TRICHLOROBENZENE	50.00	1.320	1.346	-2.0	102	0.00	16.65
130 M	1,2,4-trichlorobenzene	50.00	1.057	1.046	1.0	96	0.00	17.33
131 M	hexachlorobutadiene	50.00	0.643	0.568	11.7	96	0.00	17.44
132 M	naphthalene	50.00	1.858	1.964	-5.7	101	0.00	17.62
133 M	1,2,3-trichlorobenzene	50.00	0.892	0.907	-1.7	100	0.00	17.88
134 m	hexachloroethane	50.00	0.610	0.592	3.0	104	0.00	15.89

(#) = Out of Range
4B54249.D M4B2289.M

SPCC's out = 0 CCC's out = 0
Mon Aug 10 08:22:28 2015 GCMS4B

6.92
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Continuing Calibration Summary

Job Number: JC1107

Sample: V4B2296-CC2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54407.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Evaluate Continuing Calibration Report

Data File : C:\MSDCHEM\1\DATA\4B54407.D Vial: 3
 Acq On : 12 Aug 2015 10:09 am Operator: TOANP
 Sample : CC2289-20 Inst : MS4B
 Misc : MS89342,V4B2296,w,,,1 Multiplr: 1.00
 MS Integration Params: RTEINT.P

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 Last Update : Mon Aug 10 08:19:32 2015
 Response via : Multiple Level Calibration

Min. RRF : 0.010 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	TrueValue	AvgRF	CCRF	%Dev	Area%	Dev (min)	R.T.
1	tert butyl alcohol-d9	500.00	1.000	1.000	0.0	94	0.00	6.58
2 M	tertiary butyl alcohol	100.00	1.209	1.151	4.8	87	0.00	6.68
3 M	1,4-dioxane	500.00	0.112	0.128	-14.3	100	0.00	10.11
4	Ethanol		-----NA-----					
5 I	pentafluorobenzene	50.00	1.000	1.000	0.0	99	0.00	8.56
6	CHLOROTRIFLUOROETHENE		-----NA-----					
7 M	chlorodifluoromethane	20.00	0.506	0.451	10.9	92	0.00	3.63
8 M	dichlorodifluoromethane	20.00	0.484	0.425	12.2	92	0.00	3.61
9	Freon 114		-----NA-----					
10 M	chloromethane	20.00	0.185	0.173	6.5	98	0.01	3.97
11 M	vinyl chloride	20.00	0.532	0.458	13.9	88	0.00	4.17
12 M	bromomethane	20.00	0.305	0.319	-4.6	110	0.01	4.76
13 M	chloroethane	20.00	0.209	0.219	-4.8	106	0.01	4.92
14	vinyl bromide		-----NA-----					
15 M	trichlorofluoromethane	20.00	0.503	0.501	0.4	101	0.00	5.28
16	1,3-butadiene		-----NA-----					
17	Pentane		-----NA-----					
18	freon 123a		-----NA-----					
19 M	ethyl ether	20.00	0.191	0.163	14.7	86	0.00	5.63
20	2-chloropropane	20.00	0.061	0.056	8.2	84	0.00	5.81
21 M	acrolein	200.00	0.064	0.053	17.2	84	0.00	5.86
22 M	1,1-dichloroethene	20.00	0.395	0.374	5.3	98	0.01	6.03
23 M	acetone	20.00	0.033	0.029	12.1	86	0.00	6.05
24 M	allyl chloride	20.00	0.289	0.267	7.6	103	0.00	6.46
25 M	acetonitrile	200.00	0.032	0.027	15.6	80	0.00	6.45
26 M	iodomethane	20.00	0.746	0.695	6.8	95	0.00	6.28
27 M	carbon disulfide	20.00	1.159	1.104	4.7	95	0.00	6.38
28 M	methylene chloride	20.00	0.418	0.390	6.7	95	0.01	6.65
29 M	methyl acetate	20.00	0.056	0.051	8.9	85	0.00	6.44
30	1-chloropropane	20.00	0.714	0.616	13.7	92	0.00	6.66
31 M	methyl tert butyl ether	20.00	1.053	0.968	8.1	92	0.00	6.89
32 M	trans-1,2-dichloroethene	20.00	0.388	0.340	12.4	94	0.00	6.95
33 M	di-isopropyl ether	20.00	1.381	1.141	17.4	83	0.00	7.41
34 M	2-butanone	20.00	0.044	0.036	18.2	81	0.00	8.08
35 M	1,1-dichloroethane	20.00	0.730	0.635	13.0	88	0.00	7.46
36 M	chloroprene	20.00	0.540	0.486	10.0	92	0.00	7.55
37 M	acrylonitrile	100.00	0.149	0.126	15.4	86	0.00	6.93
38 M	vinyl acetate	20.00	0.056	0.051	8.9	89	0.00	7.43
39 M	ethyl tert-butyl ether	20.00	1.227	1.103	10.1	90	0.00	7.83
40 M	ethyl acetate	20.00	0.060	0.049	18.3	85	0.00	8.08
41 M	2,2-dichloropropane	20.00	0.309	0.285	7.8	94	0.00	8.11

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Continuing Calibration Summary

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Job Number: JC1107

Sample: V4B2296-CC2289
Lab FileID: 4B54407.D

Account: UTC United Technologies Corporation
Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

42 M	cis-1,2-dichloroethene	20.00	0.422	0.380	10.0	95	0.00	8.11
43	methylacrylate	20.00	0.057	0.050	12.3	89	0.00	8.16
44 M	propionitrile	200.00	0.060	0.050	16.7	83	0.00	8.19
45 M	bromochloromethane	20.00	0.216	0.205	5.1	96	0.00	8.40
46 M	tetrahydrofuran	20.00	0.144	0.117	18.7	82	0.00	8.43
47 M	chloroform	20.00	0.437	0.401	8.2	94	0.00	8.44
48 M	T-BUTYL FORMATE	20.00	0.279	0.256	8.2	91	0.00	8.46
49 S	dibromofluoromethane (s)	50.00	0.362	0.356	1.7	96	0.00	8.63
50 S	1,2-dichloroethane-d4 (s)	50.00	0.397	0.366	7.8	90	0.00	9.02
51 M	freon 113	20.00	0.240	0.218	9.2	92	0.01	5.97
52 M	methacrylonitrile	20.00	0.265	0.210	20.8#	81	0.00	8.34
53 M	1,1,1-trichloroethane	20.00	0.462	0.431	6.7	94	0.00	8.67
54 M	cyclohexane	20.00	0.506	0.399	21.1#	82	0.00	8.74
55	iso-butyl alcohol	200.00	0.016	0.014	12.5	86	0.00	8.83
56 I	1,4-difluorobenzene	50.00	1.000	1.000	0.0	96	0.00	9.42
57 M	epichlorohydrin	100.00	0.034	0.030	11.8	90	0.00	10.64
58 M	n-butyl alcohol	1000.00	0.010	0.010#	0.0	98	0.00	9.53
59 M	carbon tetrachloride	20.00	0.400	0.360	10.0	88	0.00	8.86
60 M	1,1-dichloropropene	20.00	0.445	0.401	9.9	89	0.00	8.83
61 M	hexane	20.00	0.338	0.283	16.3	85	0.00	7.20
62	Tert Amyl alcohol				-----NA-----			
63 M	benzene	20.00	1.309	1.210	7.6	92	0.00	9.08
64 m	iso-octane	20.00	1.051	0.813	22.6#	77	0.00	9.07
65 M	tert-amyl methyl ether	20.00	0.208	0.186	10.6	87	0.00	9.09
66 M	heptane	20.00	0.232	0.201	13.4	85	0.00	9.21
67 M	isopropyl acetate	20.00	0.156	0.130	16.7	85	0.00	8.99
68 M	1,2-dichloroethane	20.00	0.431	0.414	3.9	91	0.00	9.10
69 M	trichloroethene	20.00	0.352	0.341	3.1	98	0.00	9.74
70	Tert-amyl Ethyl Ether				-----NA-----			
71	ethyl acrylate				-----NA-----			
72 M	2-nitropropane	20.00	0.141	0.122	13.5	84	0.00	10.50
73 M	2-chloroethyl vinyl ether	100.00	0.209	0.176	15.8	80	0.00	10.50
74 M	methyl methacrylate	20.00	0.084	0.079	6.0	90	0.00	9.99
75 M	1,2-dichloropropane	20.00	0.376	0.336	10.6	87	0.00	10.00
76 M	dibromomethane	20.00	0.219	0.209	4.6	92	0.00	10.16
77 M	methylcyclohexane	20.00	0.453	0.394	13.0	86	0.00	9.94
78 M	bromodichloromethane	20.00	0.471	0.452	4.0	93	0.00	10.28
79 M	cis-1,3-dichloropropene	20.00	0.610	0.543	11.0	88	0.00	10.72
80 S	toluene-d8 (s)	50.00	1.177	1.159	1.5	94	0.00	11.02
81 M	4-methyl-2-pentanone	20.00	0.144	0.127	11.8	82	0.00	10.82
82 M	toluene	20.00	0.834	0.772	7.4	94	0.00	11.09
83 M	3-methyl-1-butanol	400.00	0.016	0.015	6.3	93	0.00	10.83
84 M	trans-1,3-dichloropropene	20.00	0.534	0.485	9.2	86	0.00	11.29
85 M	ethyl methacrylate	20.00	0.442	0.393	11.1	88	0.00	11.27
86 M	1,1,2-trichloroethane	20.00	0.290	0.265	8.6	93	0.00	11.52
87 M	2-hexanone	20.00	0.132	0.111	15.9	83	0.00	11.69
88 I	chlorobenzene-d5	50.00	1.000	1.000	0.0	96	0.00	12.60
89 M	tetrachloroethene	20.00	0.434	0.424	2.3	97	0.00	11.69
90 M	1,3-dichloropropane	20.00	0.588	0.544	7.5	90	0.00	11.71
91 M	butyl acetate	20.00	0.268	0.219	18.3	81	0.00	11.76
92 M	3,3-DIMETHYL-1-BUTANOL	200.00	0.039	0.040	-2.6	97	0.00	11.87
93 M	dibromochloromethane	20.00	0.460	0.436	5.2	95	0.00	11.99
94 M	1,2-dibromoethane	20.00	0.385	0.374	2.9	95	0.00	12.15
95	n-butyl ether				-----NA-----			
96 M	chlorobenzene	20.00	1.057	1.027	2.8	96	0.00	12.63
97 M	1,1,1,2-tetrachloroethane	20.00	0.404	0.391	3.2	96	0.00	12.69
98 M	ethylbenzene	20.00	1.761	1.695	3.7	94	0.00	12.68
99 M	m,p-xylene	40.00	0.671	0.651	3.0	95	0.00	12.80

Continuing Calibration Summary

Page 3 of 3

Job Number: JC1107

Sample: V4B2296-CC2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54407.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

100 M	o-xylene	20.00	0.677	0.676	0.1	97	0.00	13.26
101 M	styrene	20.00	1.155	1.096	5.1	93	0.00	13.27
102 M	bromoform	20.00	0.358	0.319	10.9	89	0.00	13.57
103 I	1,4-dichlorobenzene-d4	50.00	1.000	1.000	0.0	93	0.00	15.14
104 M	isopropylbenzene	20.00	3.174	3.051	3.9	95	0.00	13.63
105 S	4-bromofluorobenzene (s)	50.00	0.873	0.855	2.1	92	0.00	13.86
106	cyclohexanone	200.00	0.091	0.087	4.4	88	0.00	13.83
107 M	bromobenzene	20.00	0.931	0.932	-0.1	97	0.00	14.07
108 M	1,1,2,2-tetrachloroethane	20.00	0.930	0.890	4.3	91	0.00	13.98
109 M	trans-1,4-dichloro-2-butene	20.00	0.239	0.133	44.4#	53	0.00	14.02
110 M	1,2,3-trichloropropane	20.00	0.220	0.218	0.9	89	0.00	14.06
111 M	n-propylbenzene	20.00	3.588	3.462	3.5	94	0.00	14.08
112	4-ETHYLTOLEUENE		-----NA-----					
113 M	2-chlorotoluene	20.00	0.813	0.787	3.2	95	0.00	14.24
114 M	4-chlorotoluene	20.00	2.320	2.187	5.7	92	0.00	14.35
115 M	1,3,5-trimethylbenzene	20.00	2.672	2.593	3.0	96	0.00	14.24
116 M	tert-butylbenzene	20.00	2.318	2.315	0.1	97	0.00	14.62
117 M	pentachloroethane	20.00	0.480	0.505	-5.2	103	0.00	14.72
118 M	1,2,4-trimethylbenzene	20.00	2.574	2.542	1.2	94	0.00	14.68
119 M	sec-butylbenzene	20.00	3.522	3.378	4.1	94	0.00	14.87
120 M	1,3-dichlorobenzene	20.00	1.695	1.676	1.1	94	0.00	15.08
121 M	p-isopropyltoluene	20.00	2.944	2.908	1.2	96	0.00	15.00
122 M	1,4-dichlorobenzene	20.00	1.695	1.658	2.2	94	0.00	15.17
123	benzyl chloride	20.00	1.694	1.499	11.5	87	0.00	15.30
124 M	1,2-dichlorobenzene	20.00	1.655	1.663	-0.5	95	0.00	15.61
125	1,4-DIETHYLBENZENE		-----NA-----					
126 M	n-butylbenzene	20.00	1.460	1.385	5.1	90	0.00	15.46
127	1,2,4,5-TETRAMETHYLBENZEN		-----NA-----					
128 M	1,2-dibromo-3-chloropropane	20.00	0.149	0.137	8.1	88	0.00	16.46
129	1,3,5-TRICHLOROBENZENE	20.00	1.320	1.326	-0.5	94	0.00	16.65
130 M	1,2,4-trichlorobenzene	20.00	1.057	0.987	6.6	89	0.00	17.33
131 M	hexachlorobutadiene	20.00	0.643	0.707	-10.0	103	0.00	17.44
132 M	naphthalene	20.00	1.858	1.712	7.9	85	0.00	17.62
133 M	1,2,3-trichlorobenzene	20.00	0.892	0.847	5.0	87	0.00	17.88
134 m	hexachloroethane	20.00	0.610	0.584	4.3	94	0.00	15.89

(#) = Out of Range
4B54407.D M4B2289.M

SPCC's out = 0 CCC's out = 0
Wed Aug 12 17:18:30 2015 GCMS4B

Continuing Calibration Summary

Job Number: JC1107

Account: UTC United Technologies Corporation

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Sample: V4B2297-CC2289
Lab FileID: 4B54431.D

Page 1 of 3

Evaluate Continuing Calibration Report

Data File : C:\MSDCHEM\1\DATA\4B54431.D Vial: 27
 Acq On : 12 Aug 2015 9:35 pm Operator: TOANP
 Sample : cc2289-50 Inst : MS4B
 Misc : MS89499,V4B2297,w,,,1 Multiplr: 1.00
 MS Integration Params: RTEINT.P

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 Last Update : Mon Aug 10 08:19:32 2015
 Response via : Multiple Level Calibration

Min. RRF : 0.010 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	TrueValue	AvgRF	CCRF	*Dev	Area*	Dev (min)	R.T.
1	tert butyl alcohol-d9	500.00	1.000	1.000	0.0	91	0.00	6.58
2 M	tertiary butyl alcohol	250.00	1.209	1.305	-7.9	105	0.00	6.68
3 M	1,4-dioxane	1250.00	0.112	0.122	-8.9	98	0.00	10.11
4	Ethanol		-----NA-----					
5 I	pentafluorobenzene	50.00	1.000	1.000	0.0	97	0.00	8.56
6	CHLOROTRIFLUOROETHENE		-----NA-----					
7 M	chlorodifluoromethane	50.00	0.506	0.449	11.3	91	0.00	3.64
8 M	dichlorodifluoromethane	50.00	0.484	0.454	6.2	93	-0.01	3.60
9	Freon 114		-----NA-----					
10 M	chloromethane	50.00	0.185	0.158	14.6	87	0.00	3.96
11 M	vinyl chloride	50.00	0.532	0.463	13.0	85	0.00	4.18
12 M	bromomethane	50.00	0.305	0.296	3.0	96	0.00	4.75
13 M	chloroethane	50.00	0.209	0.197	5.7	91	0.00	4.91
14	vinyl bromide		-----NA-----					
15 M	trichlorofluoromethane	50.00	0.503	0.525	-4.4	100	0.00	5.28
16	1,3-butadiene		-----NA-----					
17	Pentane		-----NA-----					
18	freon 123a		-----NA-----					
19 M	ethyl ether	50.00	0.191	0.167	12.6	87	0.00	5.62
20	2-chloropropane	50.00	0.061	0.064	-4.9	112	0.00	5.81
21 M	acrolein	500.00	0.064	0.061	4.7	88	0.00	5.86
22 M	1,1-dichloroethene	50.00	0.395	0.356	9.9	94	0.00	6.02
23 M	acetone	50.00	0.033	0.032	3.0	93	0.00	6.05
24 M	allyl chloride	50.00	0.289	0.205	29.1#	74	0.00	6.46
25 M	acetonitrile	500.00	0.032	0.028	12.5	87	-0.01	6.44
26 M	iodomethane	50.00	0.746	0.722	3.2	96	0.00	6.27
27 M	carbon disulfide	50.00	1.159	1.097	5.3	96	0.00	6.38
28 M	methylene chloride	50.00	0.418	0.379	9.3	89	0.00	6.64
29 M	methyl acetate	50.00	0.056	0.057	-1.8	94	0.00	6.44
30	1-chloropropane	50.00	0.714	0.559	21.7#	83	0.00	6.66
31 M	methyl tert butyl ether	50.00	1.053	0.977	7.2	92	0.00	6.89
32 M	trans-1,2-dichloroethene	50.00	0.388	0.338	12.9	95	0.00	6.95
33 M	di-isopropyl ether	50.00	1.381	1.180	14.6	84	0.00	7.41
34 M	2-butanone	50.00	0.044	0.040	9.1	90	0.00	8.08
35 M	1,1-dichloroethane	50.00	0.730	0.622	14.8	89	0.00	7.45
36 M	chloroprene	50.00	0.540	0.501	7.2	98	0.00	7.55
37 M	acrylonitrile	250.00	0.149	0.130	12.8	85	0.00	6.92
38 M	vinyl acetate	50.00	0.056	0.055	1.8	96	0.00	7.43
39 M	ethyl tert-butyl ether	50.00	1.227	1.145	6.7	91	0.00	7.83
40 M	ethyl acetate	50.00	0.060	0.051	15.0	84	0.00	8.08
41 M	2,2-dichloropropane	50.00	0.309	0.266	13.9	95	0.00	8.11

Continuing Calibration Summary

Page 2 of 3

Job Number: JC1107

Sample: V4B2297-CC2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54431.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

42 M	cis-1,2-dichloroethene	50.00	0.422	0.379	10.2	92	0.00	8.11
43	methylacrylate	50.00	0.057	0.054	5.3	92	0.00	8.17
44 M	propionitrile	500.00	0.060	0.053	11.7	87	0.00	8.18
45 M	bromochloromethane	50.00	0.216	0.212	1.9	96	0.00	8.40
46 M	tetrahydrofuran	50.00	0.144	0.115	20.1#	83	0.00	8.43
47 M	chloroform	50.00	0.437	0.406	7.1	93	0.00	8.44
48 M	T-BUTYL FORMATE	50.00	0.279	0.287	-2.9	94	0.00	8.46
49 S	dibromofluoromethane (s)	50.00	0.362	0.358	1.1	96	0.00	8.63
50 S	1,2-dichloroethane-d4 (s)	50.00	0.397	0.383	3.5	94	0.00	9.02
51 M	freon 113	50.00	0.240	0.236	1.7	101	0.00	5.96
52 M	methacrylonitrile	50.00	0.265	0.218	17.7	83	0.00	8.34
53 M	1,1,1-trichloroethane	50.00	0.462	0.454	1.7	102	0.00	8.67
54 M	cyclohexane	50.00	0.506	0.427	15.6	90	0.00	8.74
55	iso-butyl alcohol	500.00	0.016	0.014	12.5	92	0.00	8.82
56 I	1,4-difluorobenzene	50.00	1.000	1.000	0.0	95	0.00	9.42
57 M	epichlorohydrin	250.00	0.034	0.032	5.9	88	0.00	10.63
58 M	n-butyl alcohol	2500.00	0.010	0.009#	10.0	99	0.00	9.53
59 M	carbon tetrachloride	50.00	0.400	0.396	1.0	101	0.00	8.86
60 M	1,1-dichloropropene	50.00	0.445	0.394	11.5	93	0.00	8.83
61 M	hexane	50.00	0.338	0.286	15.4	91	0.00	7.20
62	Tert Amyl alcohol				-----NA-----			
63 M	benzene	50.00	1.309	1.162	11.2	91	0.00	9.08
64 m	iso-octane	50.00	1.051	0.828	21.2#	82	0.00	9.06
65 M	tert-amyl methyl ether	50.00	0.208	0.208	0.0	96	0.00	9.09
66 M	heptane	50.00	0.232	0.246	-6.0	111	0.00	9.21
67 M	isopropyl acetate	50.00	0.156	0.129	17.3	88	0.00	8.99
68 M	1,2-dichloroethane	50.00	0.431	0.408	5.3	91	0.00	9.10
69 M	trichloroethene	50.00	0.352	0.331	6.0	94	0.00	9.74
70	Tert-amyl Ethyl Ether				-----NA-----			
71	ethyl acrylate				-----NA-----			
72 M	2-nitropropane	50.00	0.141	0.130	7.8	90	0.00	10.50
73 M	2-chloroethyl vinyl ether	250.00	0.209	0.195	6.7	86	0.00	10.50
74 M	methyl methacrylate	50.00	0.084	0.084	0.0	94	0.00	9.99
75 M	1,2-dichloropropane	50.00	0.376	0.328	12.8	86	0.00	10.00
76 M	dibromomethane	50.00	0.219	0.215	1.8	92	0.00	10.16
77 M	methylcyclohexane	50.00	0.453	0.436	3.8	99	0.00	9.94
78 M	bromodichloromethane	50.00	0.471	0.456	3.2	91	0.00	10.28
79 M	cis-1,3-dichloropropene	50.00	0.610	0.551	9.7	88	0.00	10.72
80 S	toluene-d8 (s)	50.00	1.177	1.165	1.0	93	0.00	11.02
81 M	4-methyl-2-pentanone	50.00	0.144	0.132	8.3	89	0.00	10.82
82 M	toluene	50.00	0.834	0.771	7.6	95	0.00	11.09
83 M	3-methyl-1-butanol	1000.00	0.016	0.014	12.5	93	0.00	10.83
84 M	trans-1,3-dichloropropene	50.00	0.534	0.508	4.9	89	0.00	11.29
85 M	ethyl methacrylate	50.00	0.442	0.428	3.2	90	0.00	11.27
86 M	1,1,2-trichloroethane	50.00	0.290	0.272	6.2	91	0.00	11.52
87 M	2-hexanone	50.00	0.132	0.115	12.9	86	0.00	11.69
88 I	chlorobenzene-d5	50.00	1.000	1.000	0.0	94	0.00	12.60
89 M	tetrachloroethene	50.00	0.434	0.419	3.5	101	0.00	11.69
90 M	1,3-dichloropropane	50.00	0.588	0.547	7.0	90	0.00	11.71
91 M	butyl acetate	50.00	0.268	0.236	11.9	87	0.00	11.76
92 M	3,3-DIMETHYL-1-BUTANOL	500.00	0.039	0.034	12.8	90	0.00	11.87
93 M	dibromochloromethane	50.00	0.460	0.450	2.2	96	0.00	11.99
94 M	1,2-dibromoethane	50.00	0.385	0.383	0.5	95	0.00	12.15
95	n-butyl ether				-----NA-----			
96 M	chlorobenzene	50.00	1.057	1.014	4.1	96	0.00	12.63
97 M	1,1,1,2-tetrachloroethane	50.00	0.404	0.390	3.5	96	0.00	12.70
98 M	ethylbenzene	50.00	1.761	1.665	5.5	95	0.00	12.69
99 M	m,p-xylene	100.00	0.671	0.638	4.9	94	0.00	12.80

Continuing Calibration Summary

Page 3 of 3

Job Number: JC1107

Sample: V4B2297-CC2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54431.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

100 M	<i>o</i> -xylene	50.00	0.677	0.669	1.2	95	0.00	13.25
101 M	styrene	50.00	1.155	1.112	3.7	94	0.00	13.27
102 M	bromoform	50.00	0.358	0.349	2.5	93	0.00	13.57
103 I	1,4-dichlorobenzene-d4	50.00	1.000	1.000	0.0	93	0.00	15.14
104 M	isopropylbenzene	50.00	3.174	2.899	8.7	95	0.00	13.63
105 S	4-bromofluorobenzene (s)	50.00	0.873	0.847	3.0	93	0.00	13.86
106	cyclohexanone	500.00	0.091	0.037	59.3#	45#	0.00	13.83
107 M	bromobenzene	50.00	0.931	0.894	4.0	95	0.00	14.07
108 M	1,1,2,2-tetrachloroethane	50.00	0.930	0.869	6.6	91	0.00	13.98
109 M	trans-1,4-dichloro-2-bute	50.00	0.239	0.148	38.1#	60	0.00	14.02
110 M	1,2,3-trichloropropane	50.00	0.220	0.215	2.3	94	0.00	14.06
111 M	<i>n</i> -propylbenzene	50.00	3.588	3.278	8.6	94	0.00	14.08
112	4-ETHYLTOLEUENE		-----NA-----					
113 M	2-chlorotoluene	50.00	0.813	0.745	8.4	94	0.00	14.24
114 M	4-chlorotoluene	50.00	2.320	2.111	9.0	91	0.00	14.35
115 M	1,3,5-trimethylbenzene	50.00	2.672	2.474	7.4	95	0.00	14.24
116 M	tert-butylbenzene	50.00	2.318	2.230	3.8	97	0.00	14.62
117 M	pentachloroethane	50.00	0.480	0.472	1.7	98	0.00	14.72
118 M	1,2,4-trimethylbenzene	50.00	2.574	2.452	4.7	94	0.00	14.68
119 M	sec-butylbenzene	50.00	3.522	3.288	6.6	95	0.00	14.87
120 M	1,3-dichlorobenzene	50.00	1.695	1.670	1.5	95	0.00	15.08
121 M	p-isopropyltoluene	50.00	2.944	2.815	4.4	96	0.00	15.00
122 M	1,4-dichlorobenzene	50.00	1.695	1.630	3.8	94	0.00	15.17
123	benzyl chloride	50.00	1.694	1.462	13.7	84	0.00	15.30
124 M	1,2-dichlorobenzene	50.00	1.655	1.652	0.2	96	0.00	15.61
125	1,4-DIETHYLBENZENE		-----NA-----					
126 M	<i>n</i> -butylbenzene	50.00	1.460	1.375	5.8	92	0.00	15.46
127	1,2,4,5-TETRAMETHYLBENZEN		-----NA-----					
128 M	1,2-dibromo-3-chloropropane	50.00	0.149	0.141	5.4	94	0.00	16.46
129	1,3,5-TRICHLOROBENZENE	50.00	1.320	1.378	-4.4	98	0.00	16.65
130 M	1,2,4-trichlorobenzene	50.00	1.057	1.071	-1.3	94	0.00	17.33
131 M	hexachlorobutadiene	50.00	0.643	0.666	-3.6	107	0.00	17.44
132 M	naphthalene	50.00	1.858	1.871	-0.7	91	0.00	17.62
133 M	1,2,3-trichlorobenzene	50.00	0.892	0.920	-3.1	96	0.00	17.88
134 m	hexachloroethane	50.00	0.610	0.576	5.6	95	0.00	15.89

(#) = Out of Range
4B54249.D M4B2289.M

SPCC's out = 0 CCC's out = 0
Thu Aug 13 12:17:21 2015 GCMS4B

Continuing Calibration Summary

Page 1 of 3

Job Number: JC1107

Sample: V4B2298-CC2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54455.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

Evaluate Continuing Calibration Report

Data File : C:\MSDCHEM\1\DATA\4B54455.D Vial: 2
 Acq On : 13 Aug 2015 9:46 am Operator: TOANP
 Sample : cc2289-20 Inst : MS4B
 Misc : MS89470,V4B2298,w,,,1 Multiplr: 1.00
 MS Integration Params: RTEINT.P

Method : C:\MSDCHEM\1\METHODS\4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 Last Update : Mon Aug 10 08:19:32 2015
 Response via : Multiple Level Calibration

Min. RRF : 0.010 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 200%

	Compound	TrueValue	AvgRF	CCRF	%Dev	Area%	Dev(min)	R.T.
1	tert butyl alcohol-d9	500.00	1.000	1.000	0.0	100	0.00	6.57
2 M	tertiary butyl alcohol	100.00	1.209	1.311	-8.4	105	0.00	6.67
3 M	1,4-dioxane	500.00	0.112	0.113	-0.9	94	-0.01	10.10
4	Ethanol		-----NA-----					
5 I	pentafluorobenzene	50.00	1.000	1.000	0.0	95	0.00	8.56
6	CHLOROTRIFLUOROETHENE		-----NA-----					
7 M	chlorodifluoromethane	20.00	0.506	0.458	9.5	90	0.00	3.64
8 M	dichlorodifluoromethane	20.00	0.484	0.562	-16.1	117	-0.01	3.60
9	Freon 114		-----NA-----					
10 M	chloromethane	20.00	0.185	0.180	2.7	99	0.01	3.97
11 M	vinyl chloride	20.00	0.532	0.474	10.9	87	0.00	4.17
12 M	bromomethane	20.00	0.305	0.344	-12.8	114	0.02	4.77
13 M	chloroethane	20.00	0.209	0.231	-10.5	107	0.02	4.92
14	vinyl bromide		-----NA-----					
15 M	trichlorofluoromethane	20.00	0.503	0.641	-27.4#	124	0.00	5.28
16	1,3-butadiene		-----NA-----					
17	Pentane		-----NA-----					
18	freon 123a		-----NA-----					
19 M	ethyl ether	20.00	0.191	0.179	6.3	91	0.00	5.63
20	2-chloropropane	20.00	0.061	0.051	16.4	74	0.00	5.81
21 M	acrolein	200.00	0.064	0.059	7.8	90	0.00	5.86
22 M	1,1-dichloroethene	20.00	0.395	0.402	-1.8	101	0.01	6.03
23 M	acetone	20.00	0.033	0.036	-9.1	104	0.00	6.05
24 M	allyl chloride	20.00	0.289	0.317	-9.7	117	0.00	6.46
25 M	acetonitrile	200.00	0.032	0.034	-6.3	95	0.00	6.45
26 M	iodomethane	20.00	0.746	0.706	5.4	92	0.01	6.28
27 M	carbon disulfide	20.00	1.159	1.088	6.1	90	0.00	6.38
28 M	methylene chloride	20.00	0.418	0.428	-2.4	100	0.02	6.66
29 M	methyl acetate	20.00	0.056	0.058	-3.6	93	0.00	6.44
30	1-chloropropane	20.00	0.714	0.641	10.2	91	0.00	6.65
31 M	methyl tert butyl ether	20.00	1.053	1.062	-0.9	97	0.00	6.89
32 M	trans-1,2-dichloroethene	20.00	0.388	0.358	7.7	95	0.00	6.95
33 M	di-isopropyl ether	20.00	1.381	1.192	13.7	83	0.00	7.41
34 M	2-butanone	20.00	0.044	0.044	0.0	94	0.00	8.08
35 M	1,1-dichloroethane	20.00	0.730	0.691	5.3	92	0.00	7.45
36 M	chloroprene	20.00	0.540	0.512	5.2	93	0.00	7.54
37 M	acrylonitrile	100.00	0.149	0.144	3.4	94	0.00	6.92
38 M	vinyl acetate	20.00	0.056	0.054	3.6	91	0.00	7.43
39 M	ethyl tert-butyl ether	20.00	1.227	1.157	5.7	91	0.00	7.82
40 M	ethyl acetate	20.00	0.060	0.054	10.0	89	-0.01	8.08
41 M	2,2-dichloropropane	20.00	0.309	0.293	5.2	93	0.00	8.10

Continuing Calibration Summary

Page 2 of 3

Job Number: JC1107

Sample: V4B2298-CC2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54455.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

42 M	cis-1,2-dichloroethene	20.00	0.422	0.413	2.1	99	0.00	8.11
43	methylacrylate	20.00	0.057	0.056	1.8	95	0.00	8.16
44 M	propionitrile	200.00	0.060	0.059	1.7	93	0.00	8.18
45 M	bromochloromethane	20.00	0.216	0.229	-6.0	102	0.00	8.40
46 M	tetrahydrofuran	20.00	0.144	0.131	9.0	88	-0.01	8.42
47 M	chloroform	20.00	0.437	0.441	-0.9	99	0.00	8.44
48 M	T-BUTYL FORMATE	20.00	0.279	0.269	3.6	92	0.00	8.46
49 S	dibromofluoromethane (s)	50.00	0.362	0.355	1.9	92	0.00	8.62
50 S	1,2-dichloroethane-d4 (s)	50.00	0.397	0.378	4.8	89	0.00	9.01
51 M	freon 113	20.00	0.240	0.280	-16.7	113	0.01	5.97
52 M	methacrylonitrile	20.00	0.265	0.235	11.3	87	0.00	8.34
53 M	1,1,1-trichloroethane	20.00	0.462	0.459	0.6	96	0.00	8.67
54 M	cyclohexane	20.00	0.506	0.488	3.6	97	0.00	8.74
55	iso-butyl alcohol	200.00	0.016	0.018	-12.5	110	0.00	8.82
56 I	1,4-difluorobenzene	50.00	1.000	1.000	0.0	91	0.00	9.41
57 M	epichlorohydrin	100.00	0.034	0.033	2.9	96	0.00	10.63
58 M	n-butyl alcohol	1000.00	0.010	0.010	0.0	97	0.00	9.52
59 M	carbon tetrachloride	20.00	0.400	0.420	-5.0	98	0.00	8.85
60 M	1,1-dichloropropene	20.00	0.445	0.451	-1.3	95	0.00	8.83
61 M	hexane	20.00	0.338	0.390	-15.4	112	0.00	7.20
62	Tert Amyl alcohol				-----NA-----			
63 M	benzene	20.00	1.309	1.294	1.1	94	0.00	9.08
64 M	iso-octane	20.00	1.051	1.020	2.9	92	0.00	9.06
65 M	tert-amyl methyl ether	20.00	0.208	0.207	0.5	92	0.00	9.09
66 M	heptane	20.00	0.232	0.347	-49.6#	141	0.00	9.21
67 M	isopropyl acetate	20.00	0.156	0.141	9.6	88	0.00	8.99
68 M	1,2-dichloroethane	20.00	0.431	0.464	-7.7	97	0.00	9.10
69 M	trichloroethene	20.00	0.352	0.362	-2.8	99	0.00	9.73
70	Tert-amyl Ethyl Ether				-----NA-----			
71	ethyl acrylate				-----NA-----			
72 M	2-nitropropane	20.00	0.141	0.142	-0.7	92	0.00	10.50
73 M	2-chloroethyl vinyl ether	100.00	0.209	0.186	11.0	80	0.00	10.50
74 M	methyl methacrylate	20.00	0.084	0.090	-7.1	98	0.00	9.98
75 M	1,2-dichloropropane	20.00	0.376	0.358	4.8	88	0.00	10.00
76 M	dibromomethane	20.00	0.219	0.236	-7.8	99	0.00	10.16
77 M	methylcyclohexane	20.00	0.453	0.520	-14.8	108	0.00	9.94
78 M	bromodichloromethane	20.00	0.471	0.496	-5.3	97	0.00	10.27
79 M	cis-1,3-dichloropropene	20.00	0.610	0.595	2.5	92	0.00	10.72
80 S	toluene-d8 (s)	50.00	1.177	1.157	1.7	90	0.00	11.01
81 M	4-methyl-2-pentanone	20.00	0.144	0.142	1.4	88	0.00	10.82
82 M	toluene	20.00	0.834	0.851	-2.0	99	0.00	11.09
83 M	3-methyl-1-butanol	400.00	0.016	0.016	0.0	95	0.00	10.83
84 M	trans-1,3-dichloropropene	20.00	0.534	0.538	-0.7	91	0.00	11.29
85 M	ethyl methacrylate	20.00	0.442	0.441	0.2	94	0.00	11.27
86 M	1,1,2-trichloroethane	20.00	0.290	0.293	-1.0	97	0.00	11.52
87 M	2-hexanone	20.00	0.132	0.123	6.8	87	0.00	11.69
88 I	chlorobenzene-d5	50.00	1.000	1.000	0.0	91	0.00	12.60
89 M	tetrachloroethene	20.00	0.434	0.451	-3.9	98	0.00	11.69
90 M	1,3-dichloropropane	20.00	0.588	0.614	-4.4	96	0.00	11.71
91 M	butyl acetate	20.00	0.268	0.235	12.3	82	0.00	11.76
92 M	3,3-DIMETHYL-1-BUTANOL	200.00	0.039	0.039	0.0	90	0.00	11.87
93 M	dibromochloromethane	20.00	0.460	0.495	-7.6	102	0.00	11.99
94 M	1,2-dibromoethane	20.00	0.385	0.418	-8.6	100	0.00	12.15
95	n-butyl ether				-----NA-----			
96 M	chlorobenzene	20.00	1.057	1.118	-5.8	99	0.00	12.63
97 M	1,1,1,2-tetrachloroethane	20.00	0.404	0.434	-7.4	101	0.00	12.69
98 M	ethylbenzene	20.00	1.761	1.835	-4.2	97	0.00	12.68
99 M	m,p-xylene	40.00	0.671	0.715	-6.6	99	0.00	12.80

Continuing Calibration Summary

Page 3 of 3

Job Number: JC1107

Sample: V4B2298-CC2289

Account: UTC United Technologies Corporation

Lab FileID: 4B54455.D

Project: ENSRILW: UTAS Plants 1/2 Facility, Rockford, IL

100 M	o-xylene	20.00	0.677	0.732	-8.1	99	0.00	13.25
101 M	styrene	20.00	1.155	1.195	-3.5	96	0.00	13.27
102 M	bromoform	20.00	0.358	0.373	-4.2	98	0.00	13.57
103 I	1,4-dichlorobenzene-d4	50.00	1.000	1.000	0.0	86	0.00	15.14
104 M	isopropylbenzene	20.00	3.174	3.421	-7.8	99	0.00	13.63
105 S	4-bromofluorobenzene (s)	50.00	0.873	0.877	-0.5	87	0.00	13.86
106	cyclohexanone	200.00	0.091	0.044	51.6#	41#	0.00	13.82
107 M	bromobenzene	20.00	0.931	1.046	-12.4	101	0.00	14.07
108 M	1,1,2,2-tetrachloroethane	20.00	0.930	1.012	-8.8	96	0.00	13.98
109 M	trans-1,4-dichloro-2-bute	20.00	0.239	0.181	24.3#	67	0.00	14.02
110 M	1,2,3-trichloropropane	20.00	0.220	0.261	-18.6	99	0.00	14.05
111 M	n-propylbenzene	20.00	3.588	3.733	-4.0	94	0.00	14.08
112	4-ETHYLTOLEUENE				-----NA-----			
113 M	2-chlorotoluene	20.00	0.813	0.862	-6.0	97	0.00	14.24
114 M	4-chlorotoluene	20.00	2.320	2.403	-3.6	94	0.00	14.35
115 M	1,3,5-trimethylbenzene	20.00	2.672	2.844	-6.4	97	0.00	14.24
116 M	tert-butylbenzene	20.00	2.318	2.548	-9.9	99	0.00	14.62
117 M	pentachloroethane	20.00	0.480	0.606	-26.3#	115	0.00	14.72
118 M	1,2,4-trimethylbenzene	20.00	2.574	2.783	-8.1	96	0.00	14.68
119 M	sec-butylbenzene	20.00	3.522	3.691	-4.8	95	0.00	14.87
120 M	1,3-dichlorobenzene	20.00	1.695	1.869	-10.3	97	0.00	15.08
121 M	p-isopropyltoluene	20.00	2.944	3.132	-6.4	96	0.00	15.00
122 M	1,4-dichlorobenzene	20.00	1.695	1.788	-5.5	94	0.00	15.17
123	benzyl chloride	20.00	1.694	1.740	-2.7	94	0.00	15.30
124 M	1,2-dichlorobenzene	20.00	1.655	1.838	-11.1	97	0.00	15.61
125	1,4-DIETHYLBENZENE				-----NA-----			
126 M	n-butylbenzene	20.00	1.460	1.434	1.8	87	0.00	15.46
127	1,2,4,5-TETRAMETHYLBENZEN				-----NA-----			
128 M	1,2-dibromo-3-chloropropane	20.00	0.149	0.159	-6.7	95	0.00	16.46
129	1,3,5-TRICHLOROBENZENE	20.00	1.320	1.446	-9.5	95	0.00	16.65
130 M	1,2,4-trichlorobenzene	20.00	1.057	1.058	-0.1	88	0.00	17.33
131 M	hexachlorobutadiene	20.00	0.643	0.768	-19.4	104	0.00	17.44
132 M	naphthalene	20.00	1.858	1.895	-2.0	88	0.00	17.62
133 M	1,2,3-trichlorobenzene	20.00	0.892	0.923	-3.5	88	0.00	17.88
134 m	hexachloroethane	20.00	0.610	0.683	-12.0	102	0.00	15.89

(#) = Out of Range
4B54248.D M4B2289.M

SPCC's out = 0 CCC's out = 0
Thu Aug 13 16:13:01 2015 GCMS4B



GC/MS Volatiles

Raw Data

7

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54425.D
 Acq On : 12 Aug 2015 6:46 pm
 Operator : TOANP
 Sample : jc1107-1
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Aug 13 11:53:59 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	138307	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	439595	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	488233	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	433020	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	211888	50.00	ug/L	0.00

System Monitoring Compounds						
49) dibromofluoromethane (s)	8.63	113	154755	48.69	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	97.38%
50) 1,2-dichloroethane-d4 (s)	9.02	65	160819	46.04	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	92.08%
80) toluene-d8 (s)	11.02	98	576118	50.12	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	100.24%
105) 4-bromofluorobenzene (s)	13.86	95	193369	52.27	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.54%

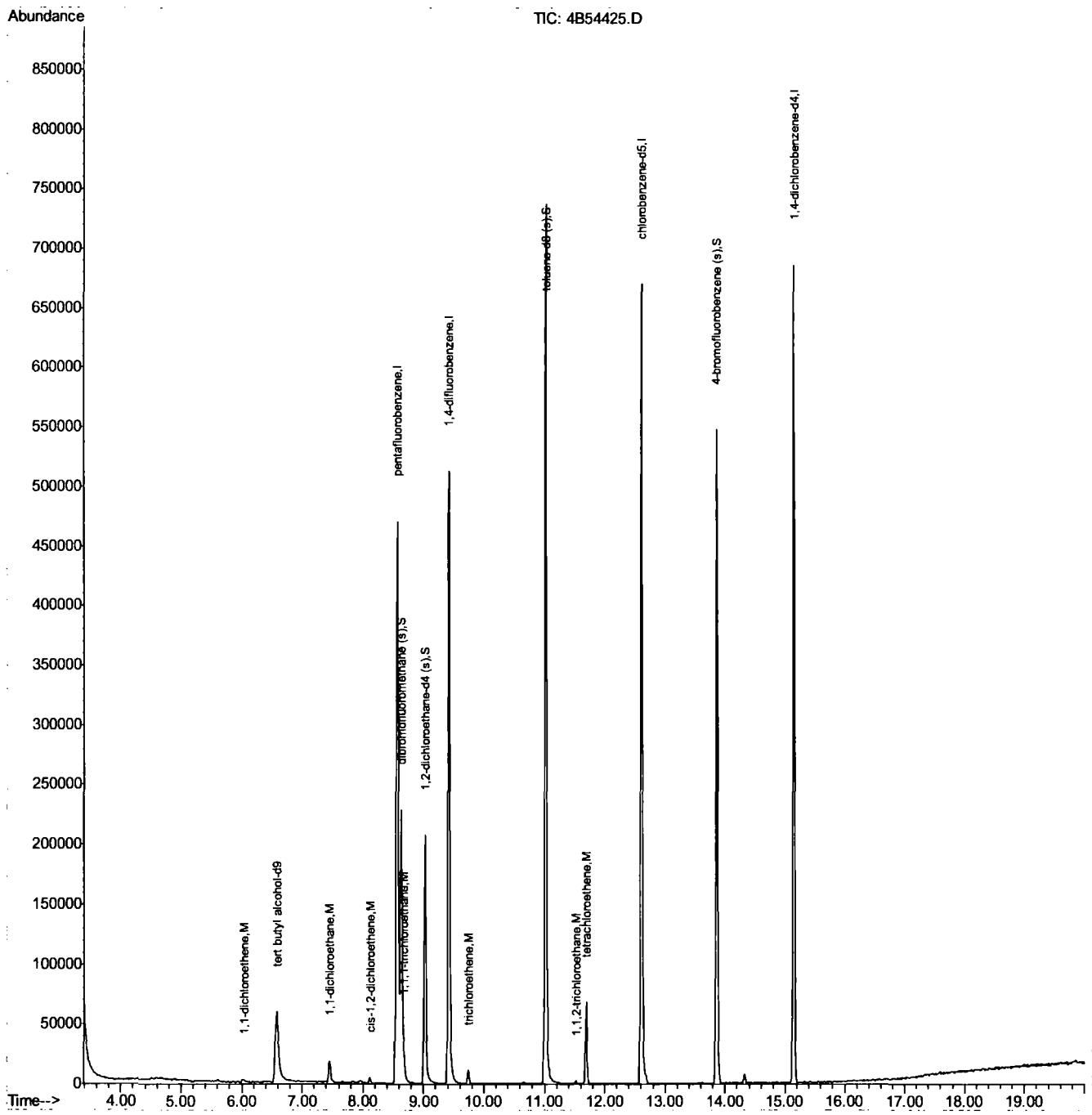
Target Compounds						Qvalue
22) 1,1-dichloroethene	6.03	96	1618	0.47	ug/L	# 47
35) 1,1-dichloroethane	7.45	63	23437	3.65	ug/L	92
42) cis-1,2-dichloroethene	8.11	96	2603	0.70	ug/L	82
53) 1,1,1-trichloroethane	8.67	97	15867	3.91	ug/L	93
69) trichloroethene	9.74	95	4280	1.24	ug/L	87
86) 1,1,2-trichloroethane	11.52	83	611	0.22	ug/L	85
89) tetrachloroethene	11.69	164	19538	5.19	ug/L	97

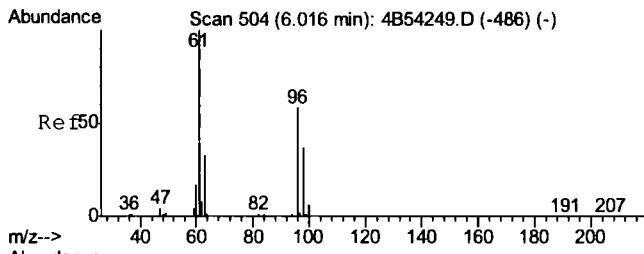
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

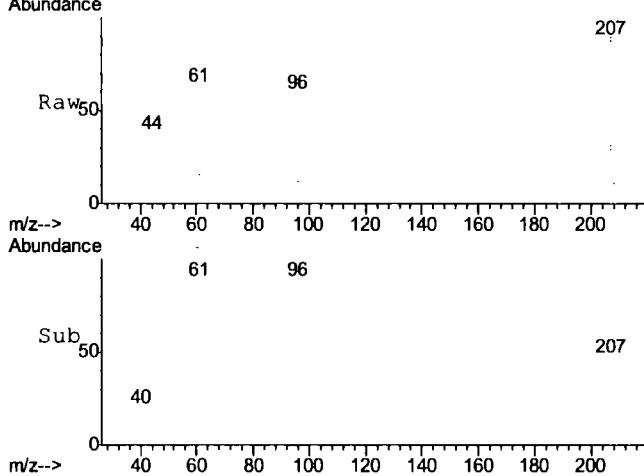
Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54425.D
 Acq On : 12 Aug 2015 6:46 pm
 Operator : TOANP
 Sample : jc1107-1
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 21 Sample Multiplier: 1

Quant Time: Aug 13 11:53:59 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

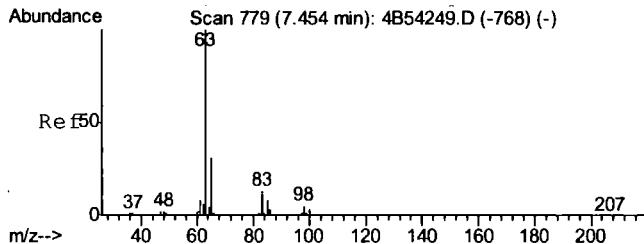




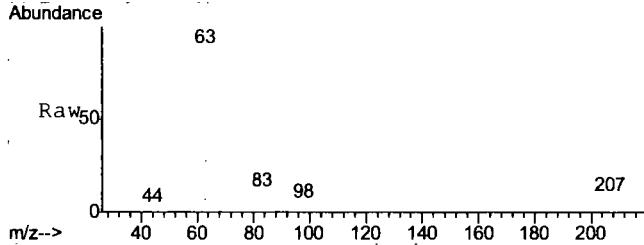
#22
1,1-dichloroethene
Concen: 0.47 ug/L
RT: 6.03 min Scan# 506
Delta R.T. 0.01 min
Lab File: 4B54425.D
Acq: 12 Aug 2015 6:46 pm



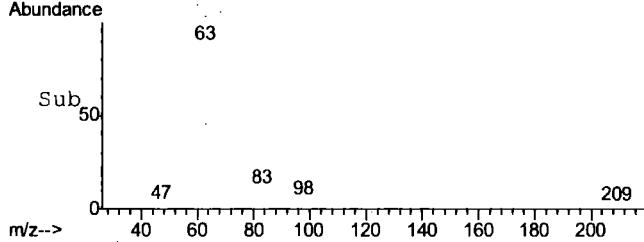
Tgt Ion: 96 Resp: 1618
Ion Ratio Lower Upper
96 100
61 84.9 141.9 201.9#
63 39.6 26.8 86.8



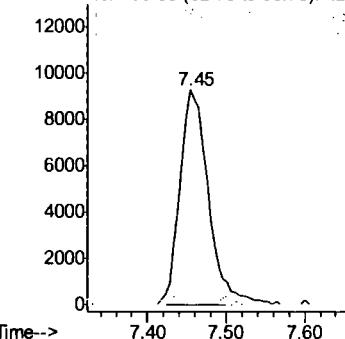
#35
1,1-dichloroethane
Concen: 3.65 ug/L
RT: 7.45 min Scan# 779
Delta R.T. -0.00 min
Lab File: 4B54425.D
Acq: 12 Aug 2015 6:46 pm

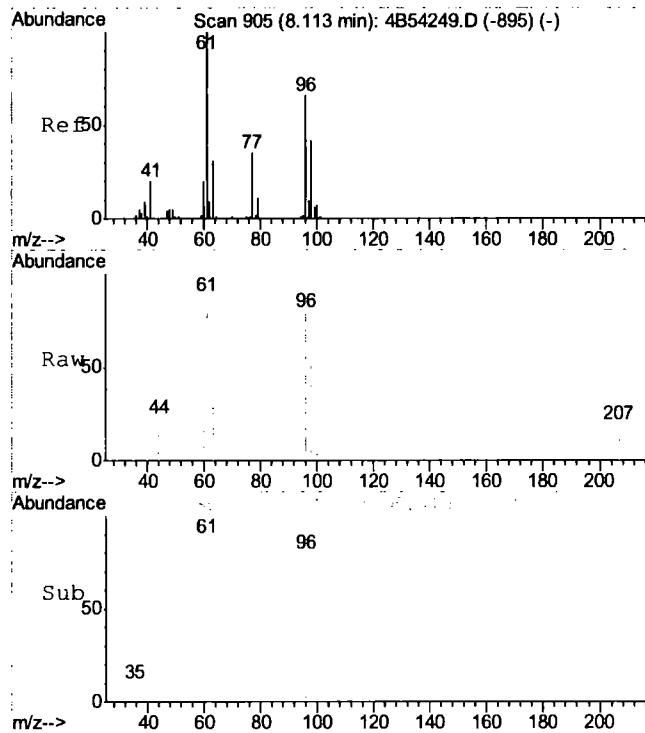


Tgt Ion: 63 Resp: 23437
Ion Ratio Lower Upper
63 100
65 36.4 1.3 61.3
83 11.2 0.0 42.6



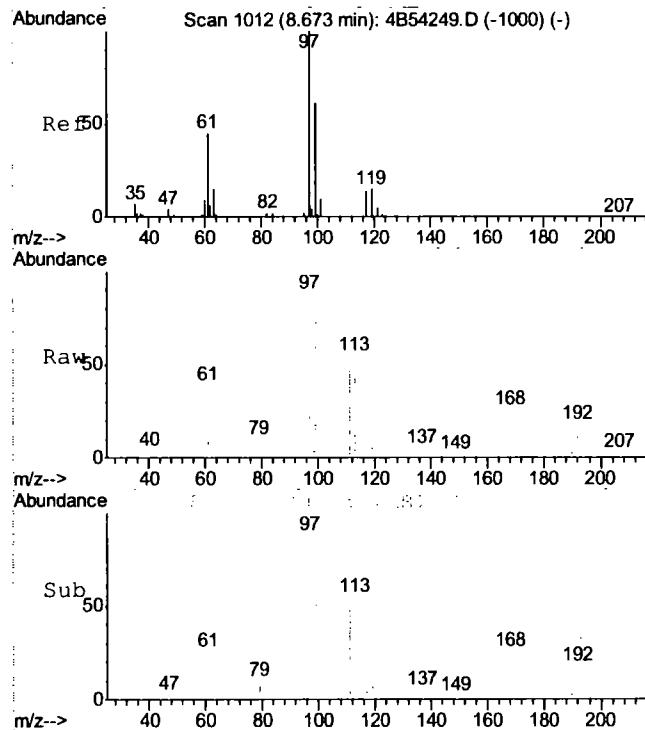
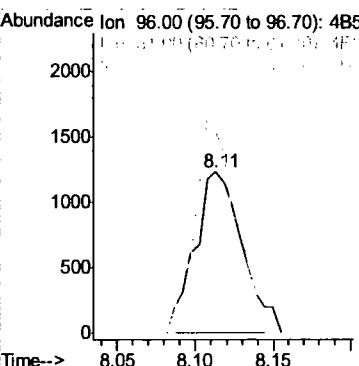
Abundance Ion 63.00 (62.70 to 63.70): 4B5





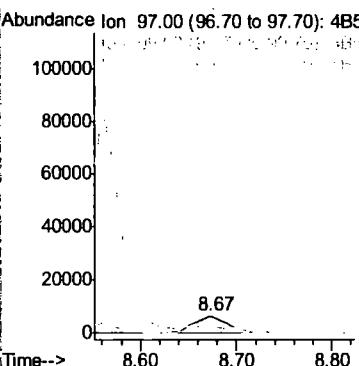
#42
cis-1,2-dichloroethene
Concen: 0.70 ug/L
RT: 8.11 min Scan# 905
Delta R.T. -0.00 min
Lab File: 4B54425.D
Acq: 12 Aug 2015 6:46 pm

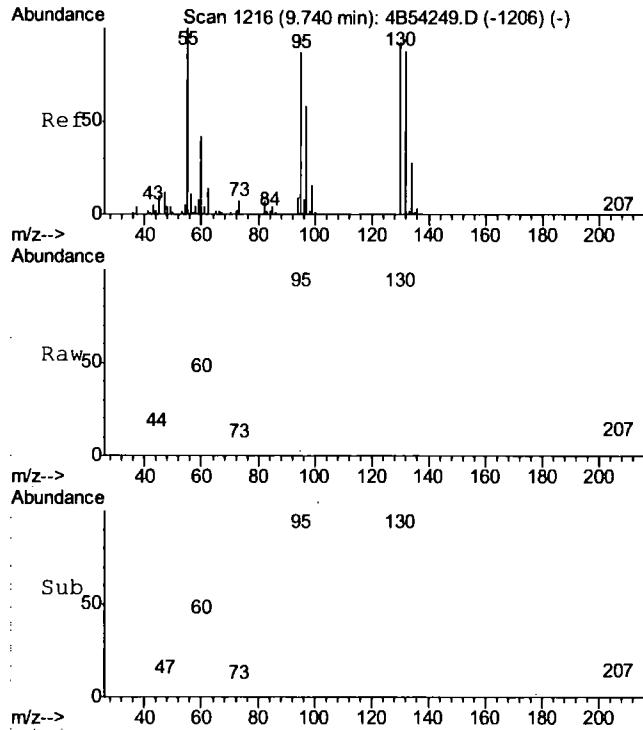
Tgt Ion: 96 Resp: 2603
Ion Ratio Lower Upper
96 100
61 124.9 123.9 183.9
98 69.3 34.0 94.0



#53
1,1,1-trichloroethane
Concen: 3.91 ug/L
RT: 8.67 min Scan# 1012
Delta R.T. -0.00 min
Lab File: 4B54425.D
Acq: 12 Aug 2015 6:46 pm

Tgt Ion: 97 Resp: 15867
Ion Ratio Lower Upper
97 100
99 70.9 35.6 95.6
61 39.4 14.6 74.6

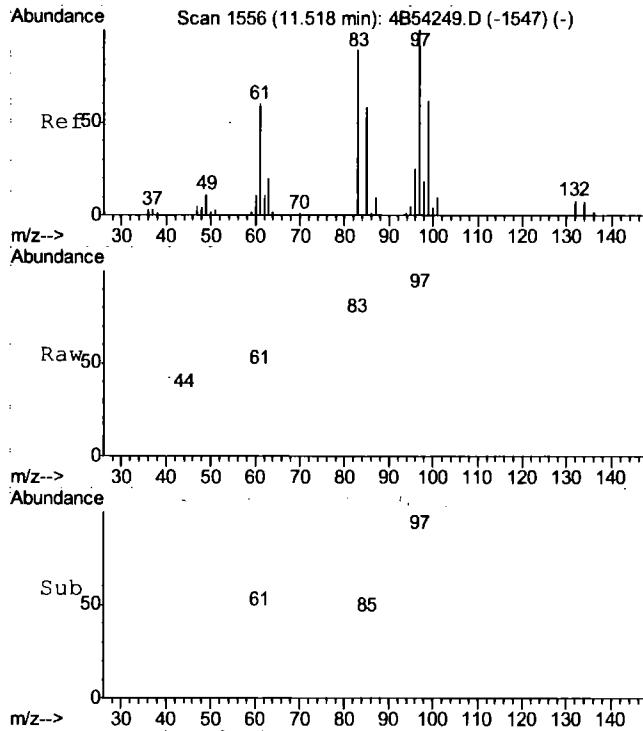
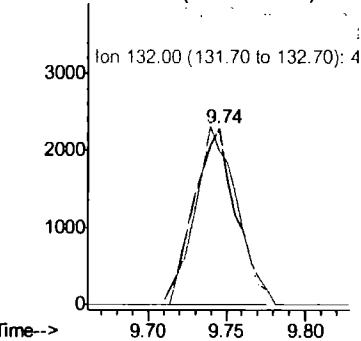




#69
trichloroethene
Concen: 1.24 ug/L
RT: 9.74 min Scan# 1217
Delta R.T. 0.01 min
Lab File: 4B54425.D
Acq: 12 Aug 2015 6:46 pm

Tgt Ion: 95 Resp: 4280
Ion Ratio Lower Upper
95 100
97 57.4 36.5 96.5
130 93.4 76.6 136.6
132 87.3 71.3 131.3

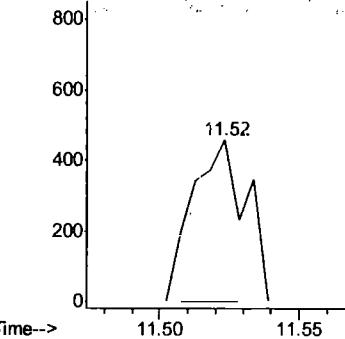
Abundance Ion 95.00 (94.70 to 95.70): 4B5

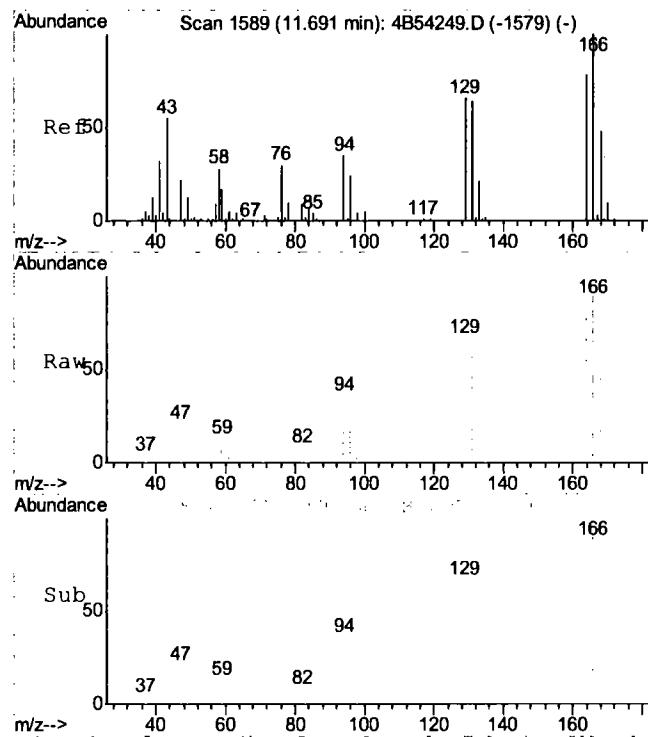


#86
1,1,2-trichloroethane
Concen: 0.22 ug/L
RT: 11.52 min Scan# 1557
Delta R.T. 0.01 min
Lab File: 4B54425.D
Acq: 12 Aug 2015 6:46 pm

Tgt Ion: 83 Resp: 611
Ion Ratio Lower Upper
83 100
97 132.8 82.1 142.1
85 58.5 34.8 94.8

Abundance Ion 83.00 (82.70 to 83.70): 4B5



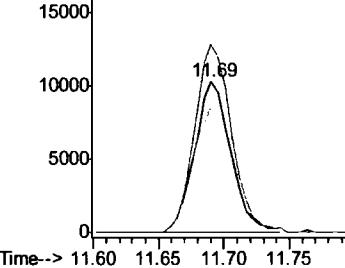


#89
tetrachloroethene
Concen: 5.19 ug/L
RT: 11.69 min Scan# 1589
Delta R.T. -0.00 min
Lab File: 4B54425.D
Acq: 12 Aug 2015 6:46 pm

Tgt Ion:164 Resp: 19538
Ion Ratio Lower Upper

164	100		
129	83.1	54.7	114.7
131	78.9	52.0	112.0
166	124.8	98.8	158.8

Abundance Ion 164.00 (163.70 to 164.70): 4
Ion 166.00 (165.70 to 166.70): 4



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54426.D
 Acq On : 12 Aug 2015 7:14 pm
 Operator : TOANP
 Sample : jc1107-2
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 22 Sample Multiplier: 1

Quant Time: Aug 13 11:54:35 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

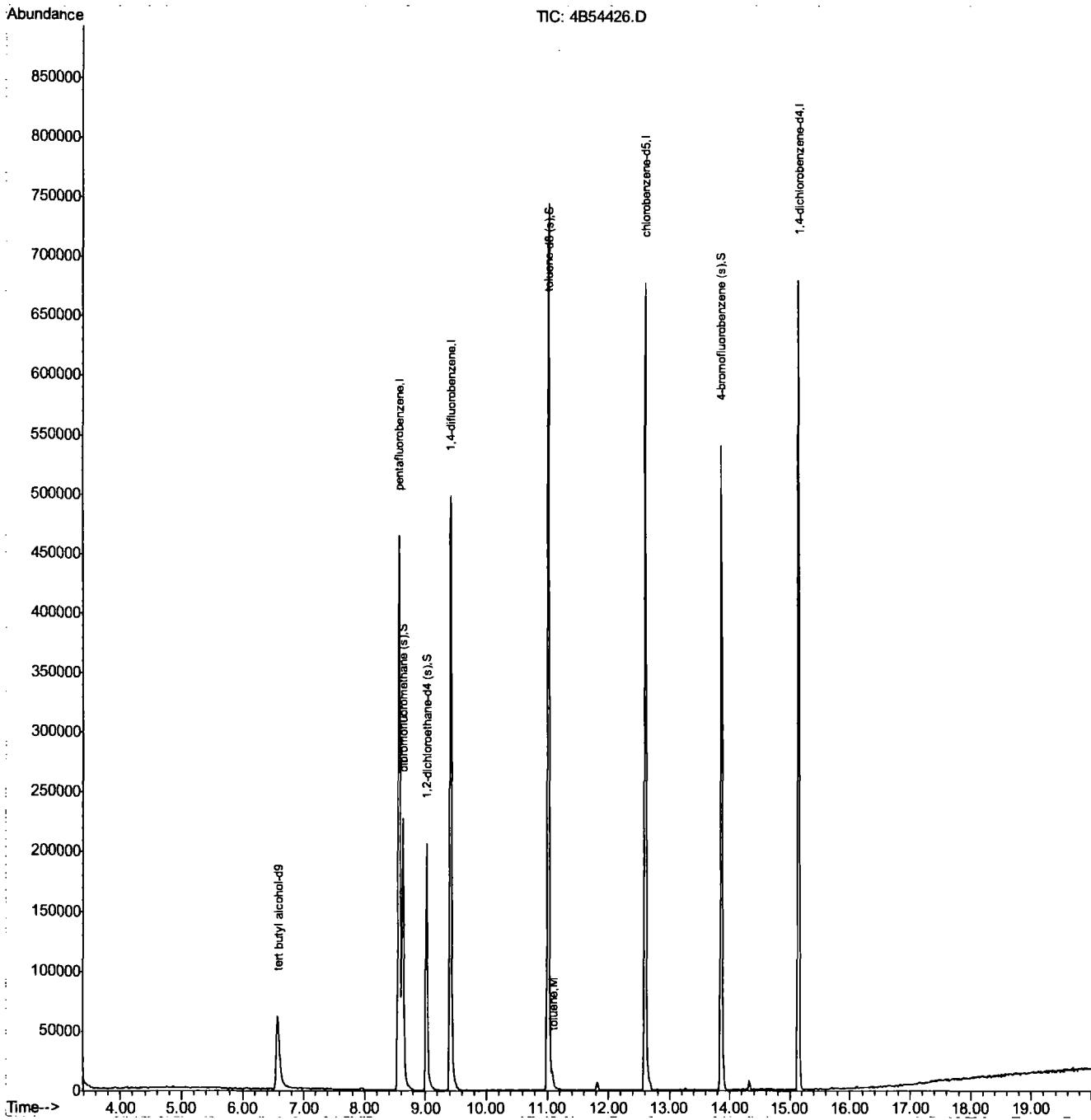
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	140461	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	437134	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	492560	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	432075	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	213576	50.00	ug/L	0.00
<hr/>						
System Monitoring Compounds						
49) dibromofluoromethane (s)	8.63	113	154048	48.74	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	97.48%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	162717	46.84	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	93.68%	
80) toluene-d8 (s)	11.02	98	574141	49.51	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	99.02%	
105) 4-bromofluorobenzene (s)	13.86	95	193164	51.80	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	103.60%	
<hr/>						
Target Compounds						
82) toluene	11.09	92	2719	0.33	ug/L	97
<hr/>						

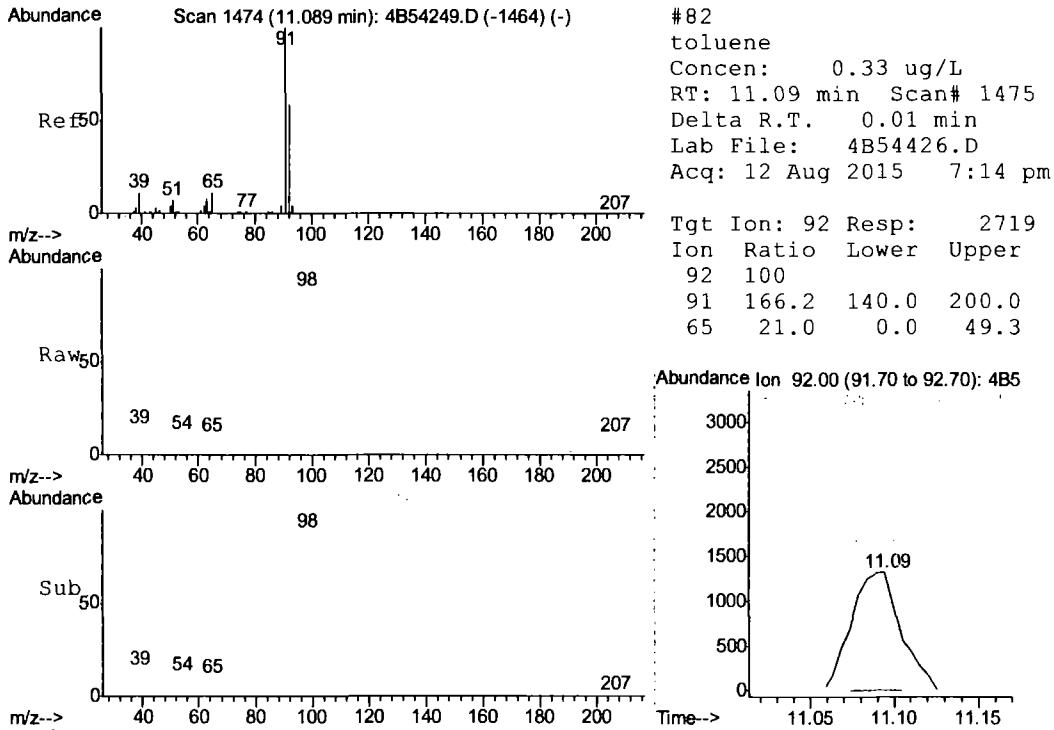
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54426.D
Acq On : 12 Aug 2015 7:14 pm
Operator : TOANP
Sample : jc1107-2
Misc : MS89470,V4B2296,w,,,1
ALS Vial : 22 Sample Multiplier: 1

Quant Time: Aug 13 11:54:35 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54416.D
 Acq On : 12 Aug 2015 2:35 pm
 Operator : TOANP
 Sample : jc1107-3
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Aug 12 17:33:24 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) tert butyl alcohol-d9	6.58	65	146811	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	443287	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	502337	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	441128	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	222699	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	157207	49.04	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	98.08%
50) 1,2-dichloroethane-d4 (s)	9.02	65	162622	46.16	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	92.32%
80) toluene-d8 (s)	11.02	98	583347	49.33	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	98.66%
105) 4-bromofluorobenzene (s)	13.86	95	202204	52.01	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.02%

Target Compounds

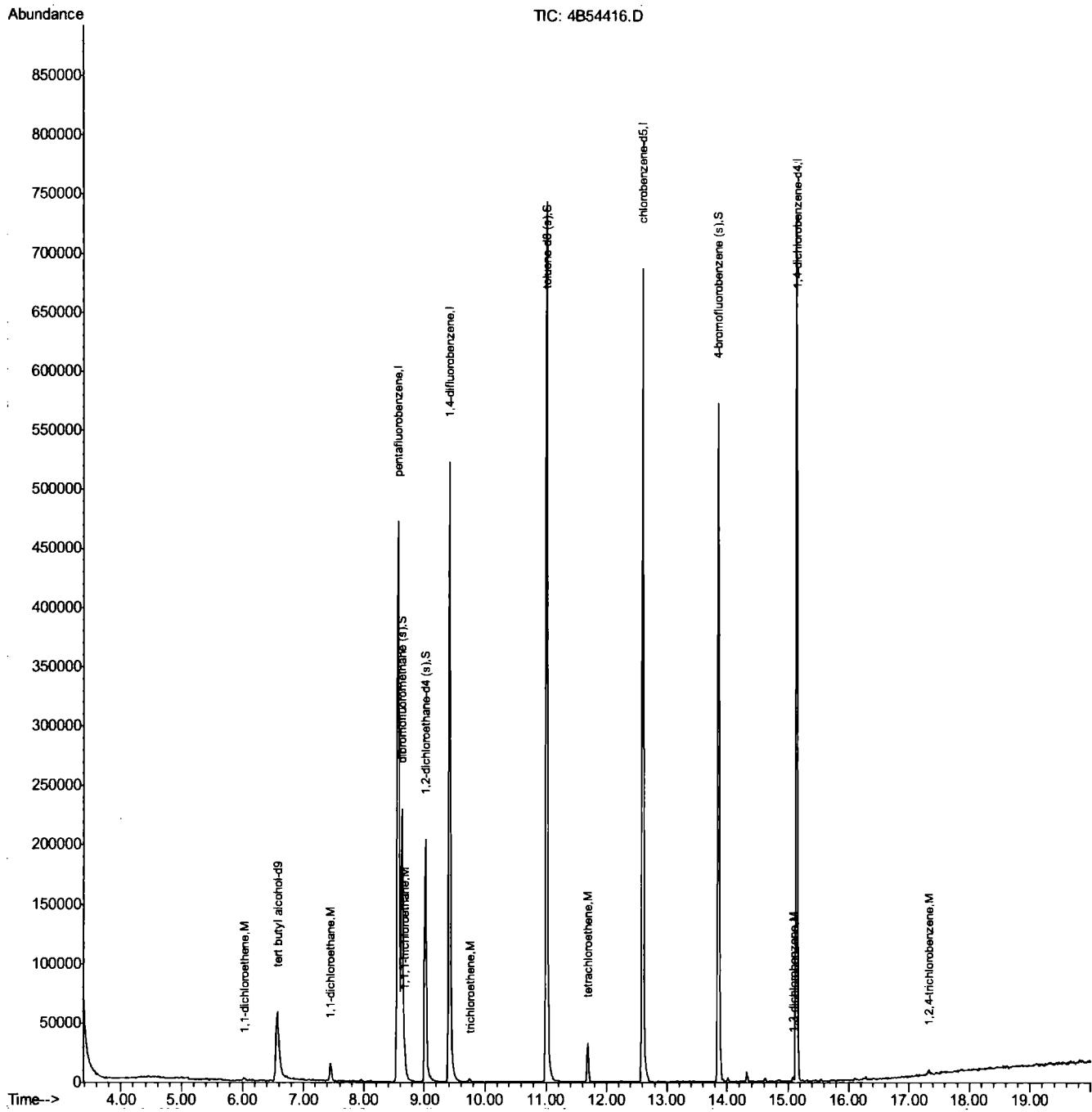
				Qvalue
22) 1,1-dichloroethene	6.03	96	1583	0.45 ug/L # 50
35) 1,1-dichloroethane	7.45	63	18549	2.87 ug/L 95
53) 1,1,1-trichloroethane	8.67	97	13114	3.20 ug/L 97
69) trichloroethene	9.74	95	1468	0.41 ug/L 87
89) tetrachloroethene	11.69	164	9638	2.51 ug/L 92
120) 1,3-dichlorobenzene	15.09	146	1820	0.24 ug/L 83
130) 1,2,4-trichlorobenzene	17.33	180	1776	0.38 ug/L 97

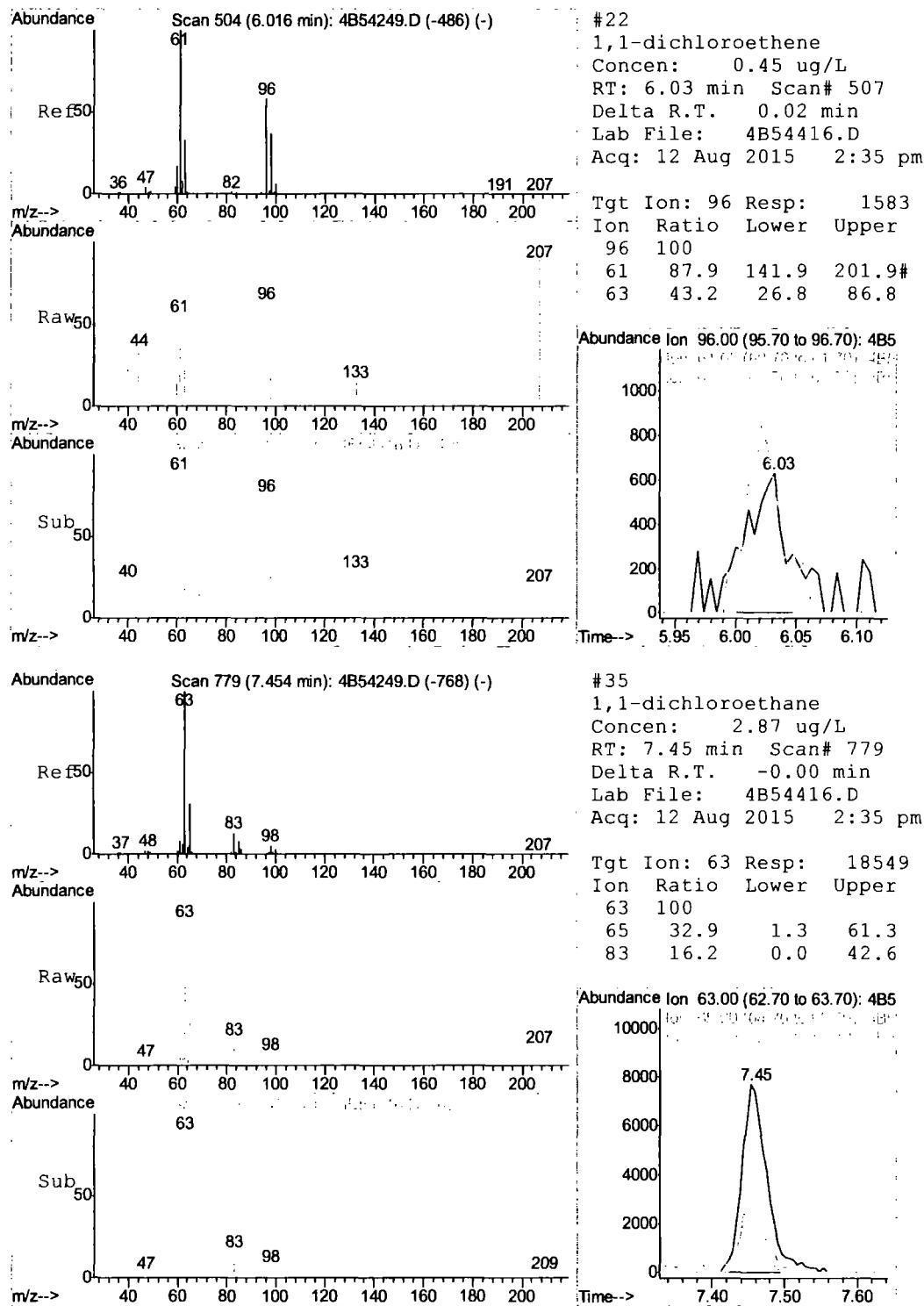
(#) = qualifier out of range (m) = manual integration (+) = signals summed

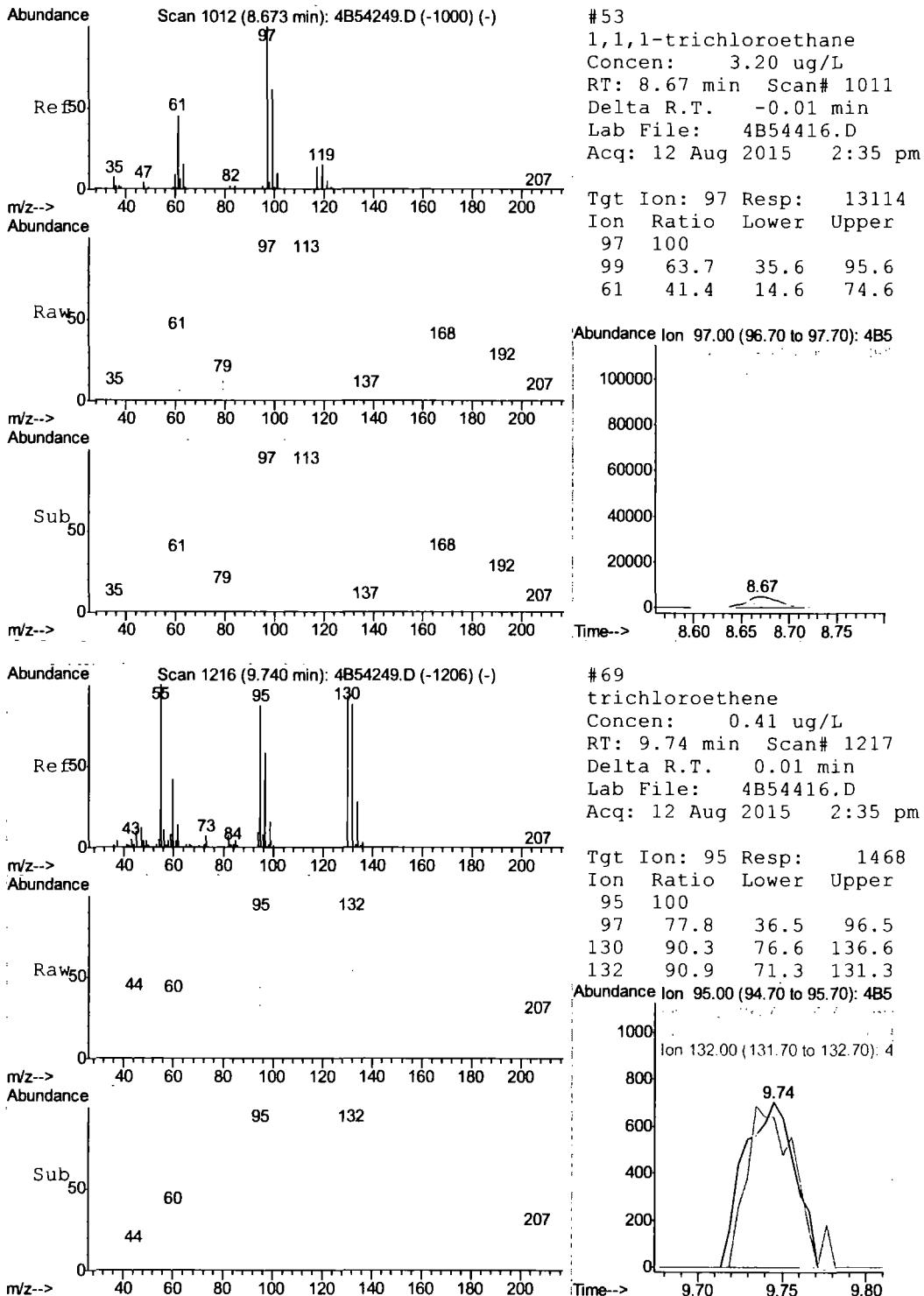
Quantitation Report (QT Reviewed)

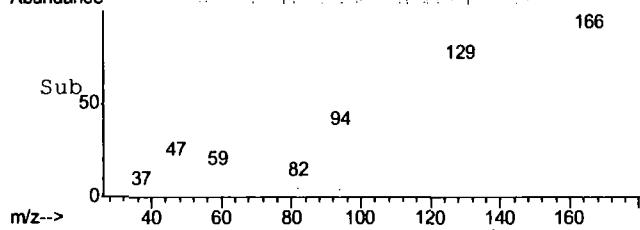
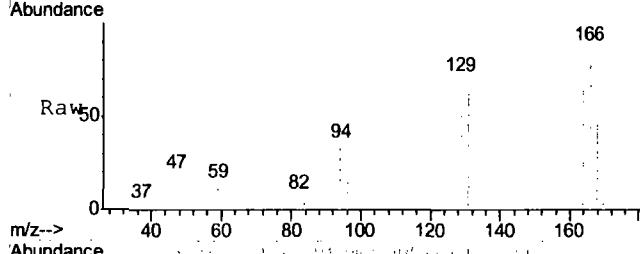
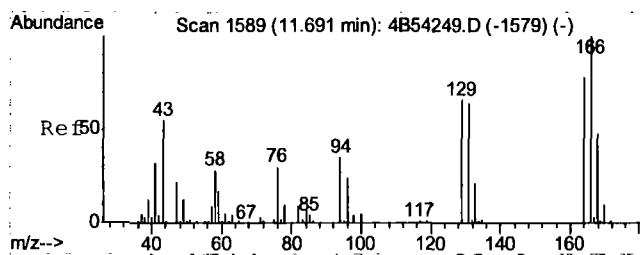
Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54416.D
 Acq On : 12 Aug 2015 2:35 pm
 Operator : TOANP
 Sample : jc1107-3
 Misc : MS89470, V4B2296, w,,,1
 ALS Vial : 12 Sample Multiplier: 1

Quant Time: Aug 12 17:33:24 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration







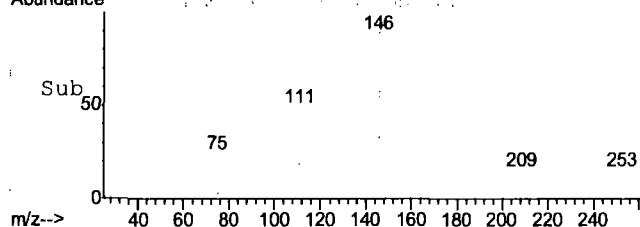
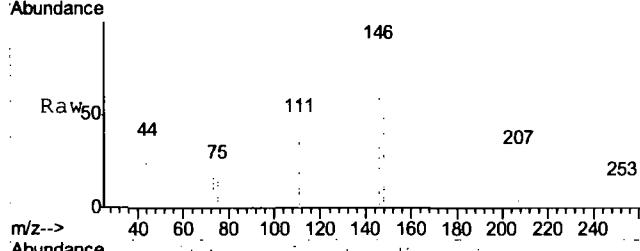
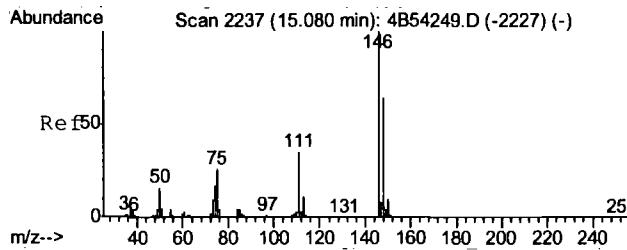
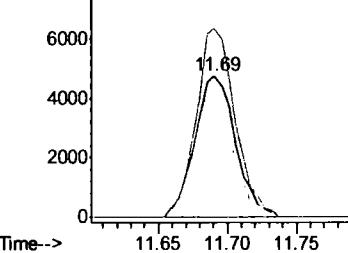


#89
tetrachloroethene
Concen: 2.51 ug/L
RT: 11.69 min Scan# 1589
Delta R.T. -0.00 min
Lab File: 4B54416.D
Acq: 12 Aug 2015 2:35 pm

Tgt Ion:164 Resp: 9638
Ion Ratio Lower Upper

164	100		
129	95.9	54.7	114.7
131	89.5	52.0	112.0
166	133.3	98.8	158.8

Abundance Ion 164.00 (163.70 to 164.70): 4
Ion 129.00 (128.70 to 130.70): 1
Ion 131.00 (130.70 to 132.70): 1
Ion 166.00 (165.70 to 166.70): 4

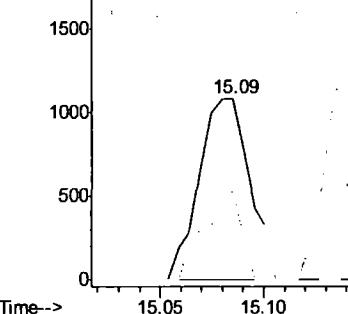


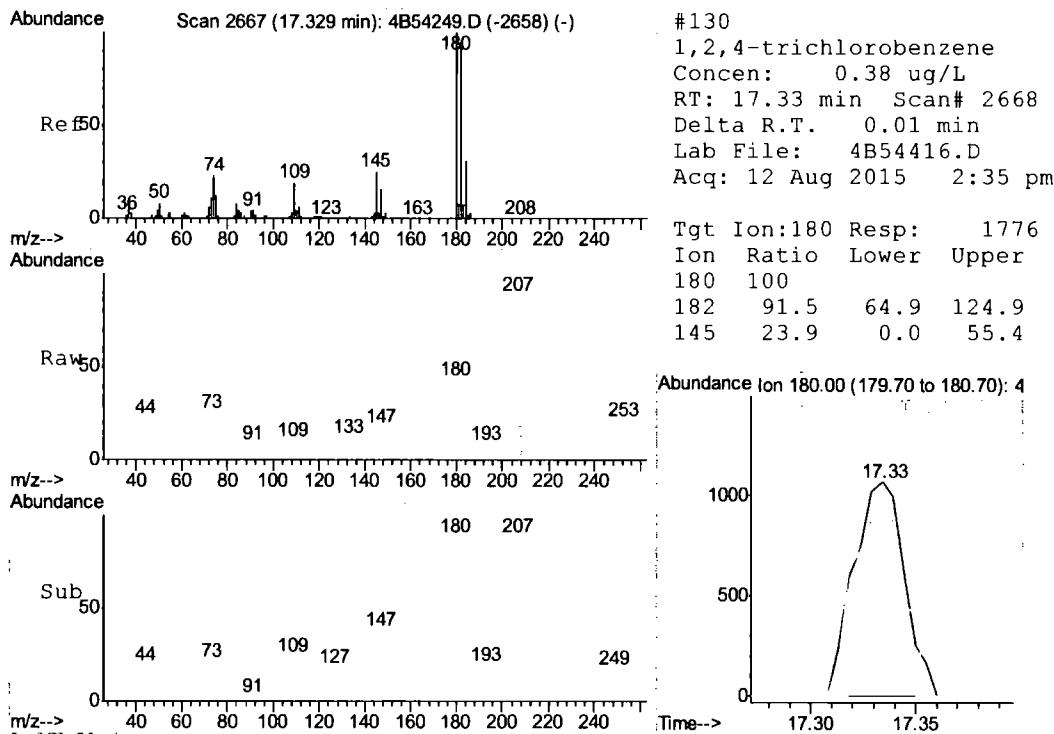
#120
1,3-dichlorobenzene
Concen: 0.24 ug/L
RT: 15.09 min Scan# 2238
Delta R.T. 0.01 min
Lab File: 4B54416.D
Acq: 12 Aug 2015 2:35 pm

Tgt Ion:146 Resp: 1820
Ion Ratio Lower Upper

146	100		
111	49.4	5.0	65.0
148	54.2	33.6	93.6

Abundance Ion 146.00 (145.70 to 146.70): 4





Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54427.D
 Acq On : 12 Aug 2015 7:42 pm
 Operator : TOANP
 Sample : jc1107-4
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 23 Sample Multiplier: 1

Quant Time: Aug 13 11:55:29 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) tert butyl alcohol-d9	6.58	65	150775	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	439086	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	491241	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	432149	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	214304	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	154067	48.52	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	97.04%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	162055	46.44	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	92.88%	
80) toluene-d8 (s)	11.02	98	571699	49.44	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	98.88%	
105) 4-bromofluorobenzene (s)	13.86	95	195212	52.18	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	104.36%	

Target Compounds

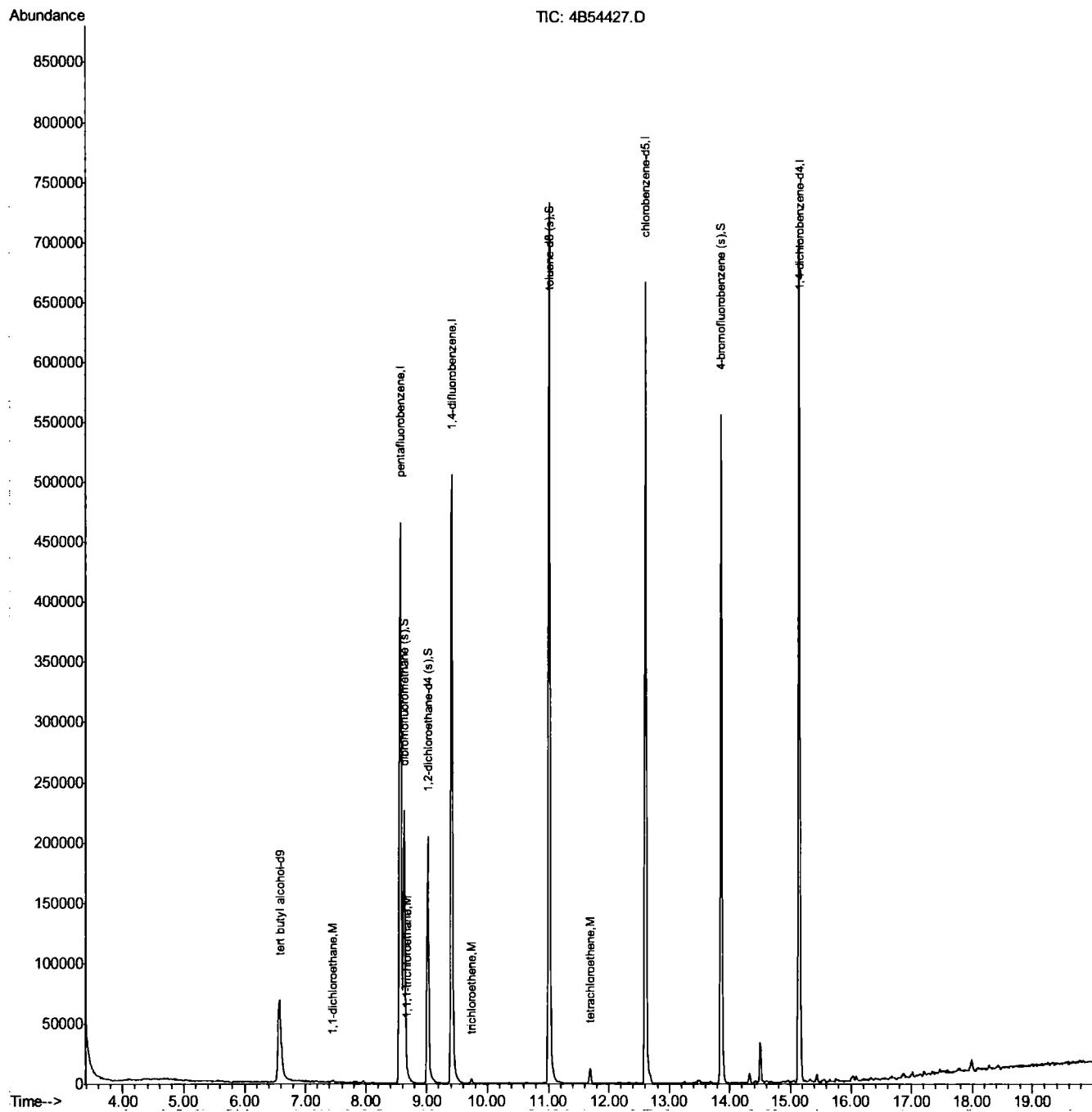
				Qvalue
35) 1,1-dichloroethane	7.45	63	2481	0.39 ug/L 95
53) 1,1,1-trichloroethane	8.68	97	1751	0.43 ug/L 95
69) trichloroethene	9.74	95	1503	0.43 ug/L 88
89) tetrachloroethene	11.70	164	4000	1.07 ug/L 87

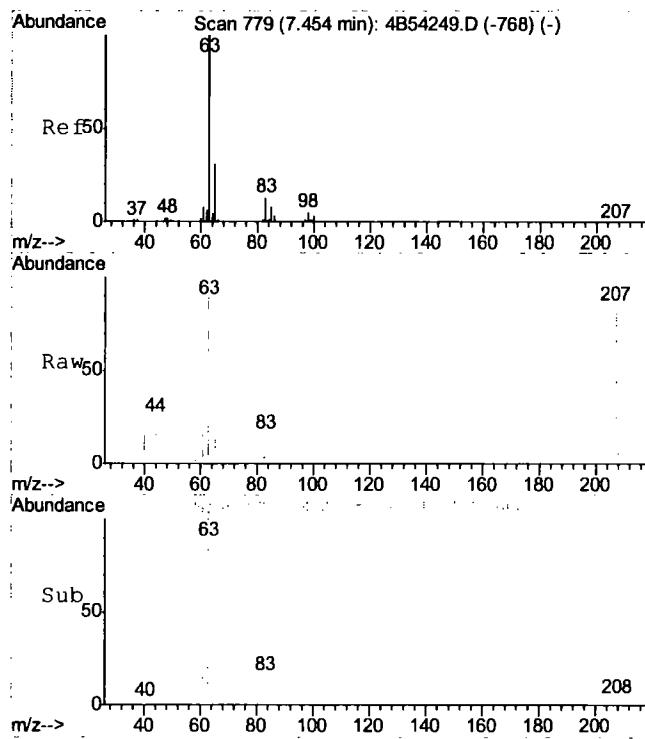
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54427.D
 Acq On : 12 Aug 2015 7:42 pm
 Operator : TOANP
 Sample : jc1107-4
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 23 Sample Multiplier: 1

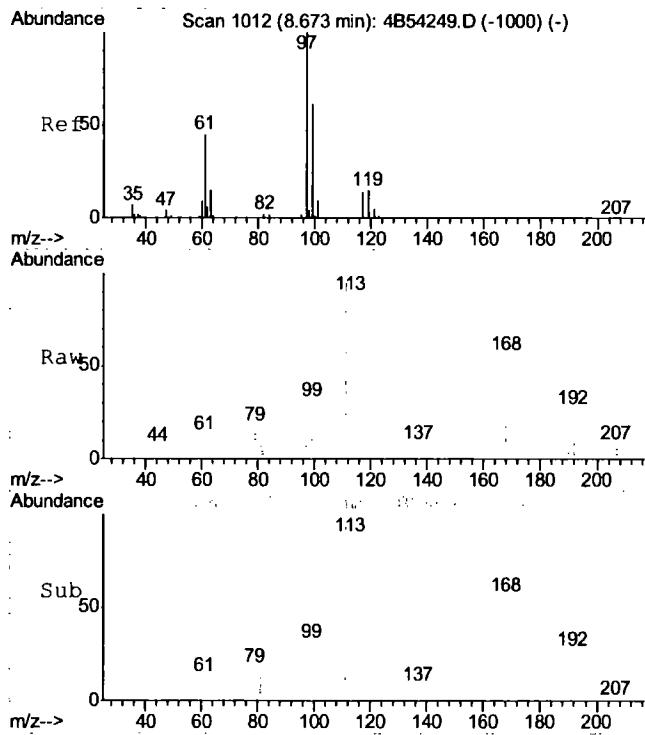
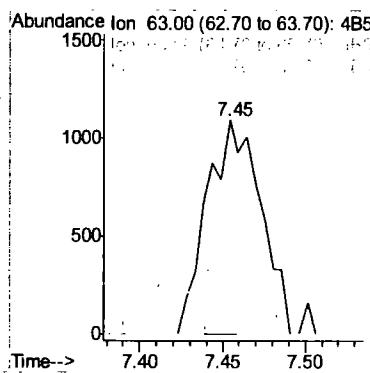
Quant Time: Aug 13 11:55:29 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration





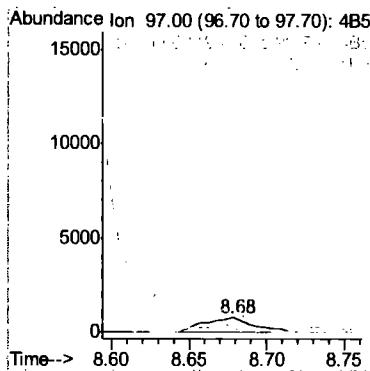
#35
1,1-dichloroethane
Concen: 0.39 ug/L
RT: 7.45 min Scan# 779
Delta R.T. -0.00 min
Lab File: 4B54427.D
Acq: 12 Aug 2015 7:42 pm

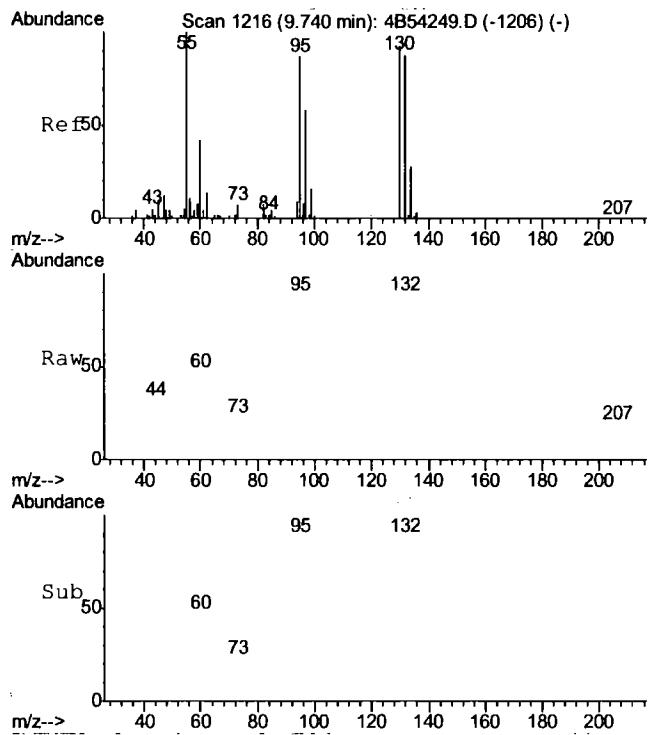
Tgt Ion: 63 Resp: 2481
Ion Ratio Lower Upper
63 100
65 32.9 1.3 61.3
83 16.2 0.0 42.6



#53
1,1,1-trichloroethane
Concen: 0.43 ug/L
RT: 8.68 min Scan# 1013
Delta R.T. 0.01 min
Lab File: 4B54427.D
Acq: 12 Aug 2015 7:42 pm

Tgt Ion: 97 Resp: 1751
Ion Ratio Lower Upper
97 100
99 59.0 35.6 95.6
61 44.9 14.6 74.6

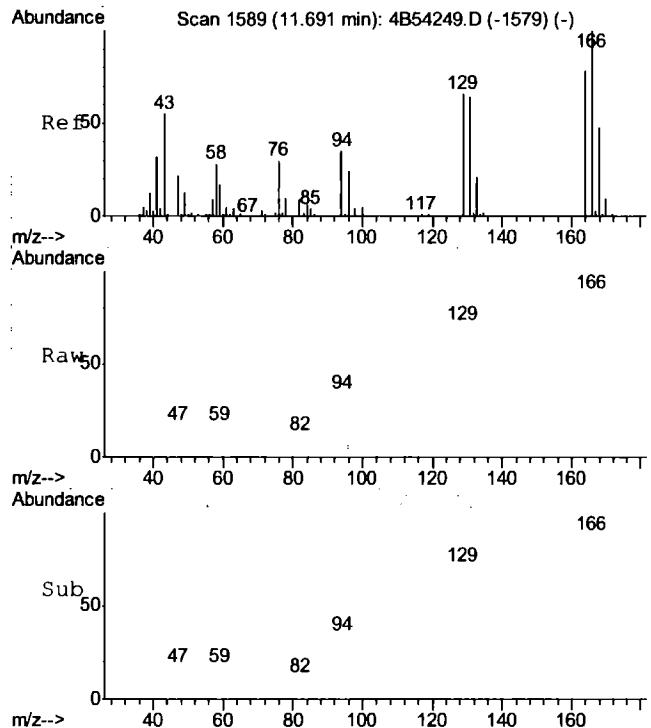
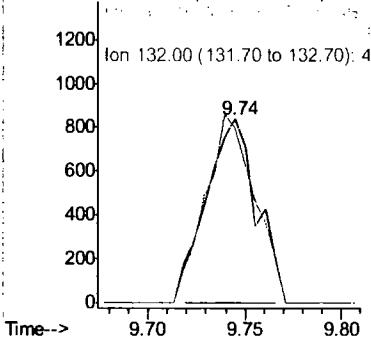




#69
trichloroethene
Concen: 0.43 ug/L
RT: 9.74 min Scan# 1217
Delta R.T. 0.01 min
Lab File: 4B54427.D
Acq: 12 Aug 2015 7:42 pm

Tgt Ion: 95 Resp: 1503
Ion Ratio Lower Upper
95 100
97 52.9 36.5 96.5
130 91.6 76.6 136.6
132 95.5 71.3 131.3

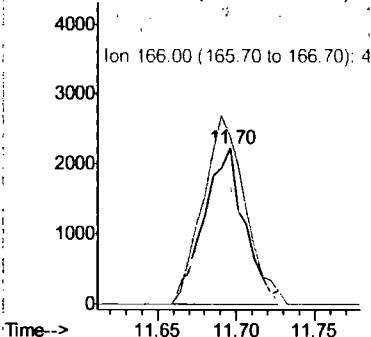
Abundance Ion 95.00 (94.70 to 95.70): 4B5



#89
tetrachloroethene
Concen: 1.07 ug/L
RT: 11.70 min Scan# 1590
Delta R.T. 0.01 min
Lab File: 4B54427.D
Acq: 12 Aug 2015 7:42 pm

Tgt Ion: 164 Resp: 4000
Ion Ratio Lower Upper
164 100
129 76.3 54.7 114.7
131 74.5 52.0 112.0
166 107.6 98.8 158.8

Abundance Ion 164.00 (163.70 to 164.70): 4



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54428.D
 Acq On : 12 Aug 2015 8:10 pm
 Operator : TOANP
 Sample : jc1107-5
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Aug 13 11:56:11 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	149773	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	427592	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	484870	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	426026	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	211603	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	151366	48.96	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	97.92%
50) 1,2-dichloroethane-d4 (s)	9.02	65	160044	47.10	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	94.20%
80) toluene-d8 (s)	11.02	98	565579	49.55	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	99.10%
105) 4-bromofluorobenzene (s)	13.86	95	192511	52.11	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.22%

Target Compounds

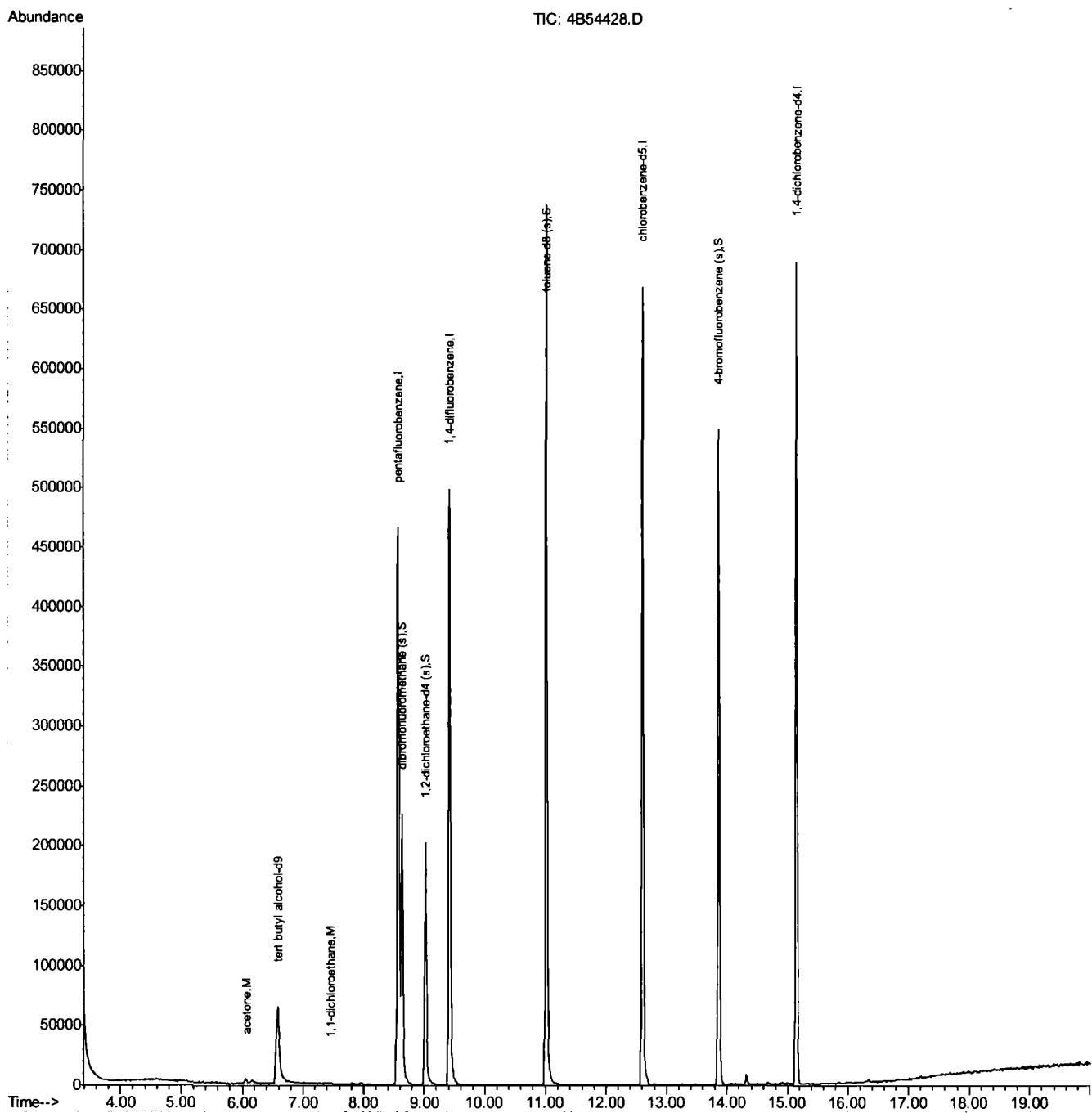
					Qvalue
23) acetone	6.06	58	2209	7.78	ug/L 97
35) 1,1-dichloroethane	7.45	63	1530	0.25	ug/L 58

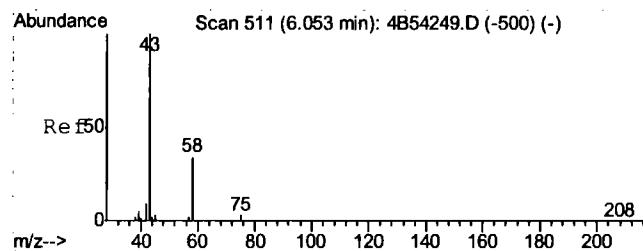
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

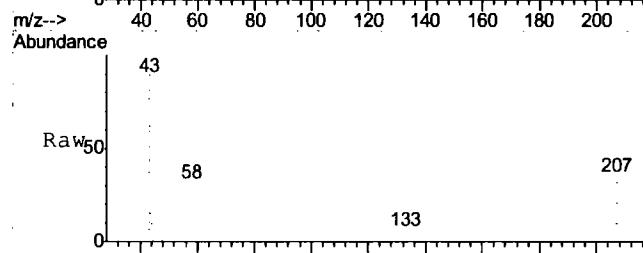
Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54428.D
 Acq On : 12 Aug 2015 8:10 pm
 Operator : TOANP
 Sample : jc1107-5
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 24 Sample Multiplier: 1

Quant Time: Aug 13 11:56:11 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

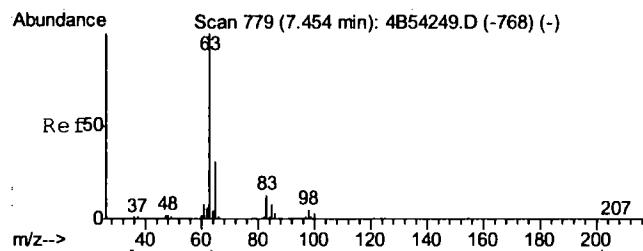
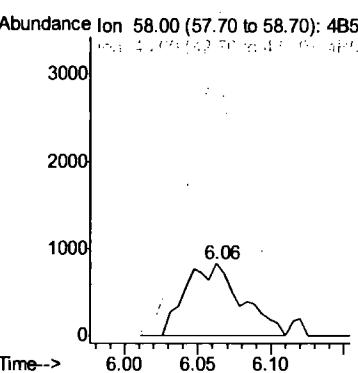
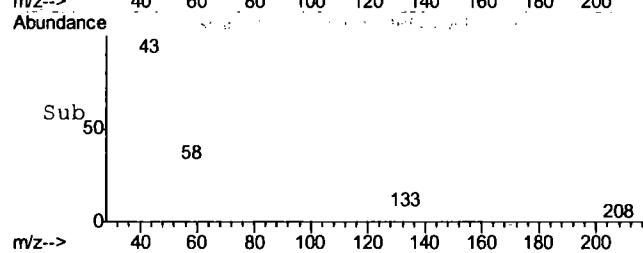




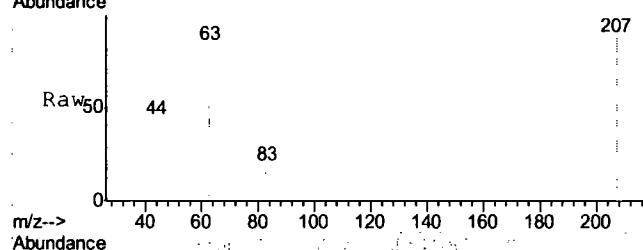
#23
acetone
Concen: 7.78 ug/L
RT: 6.06 min Scan# 513
Delta R.T. 0.01 min
Lab File: 4B54428.D
Acq: 12 Aug 2015 8:10 pm



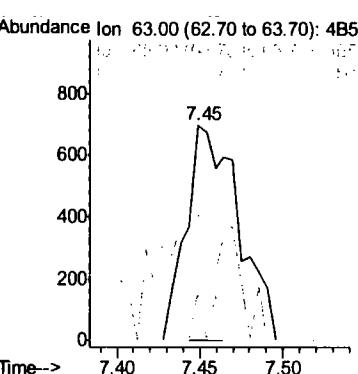
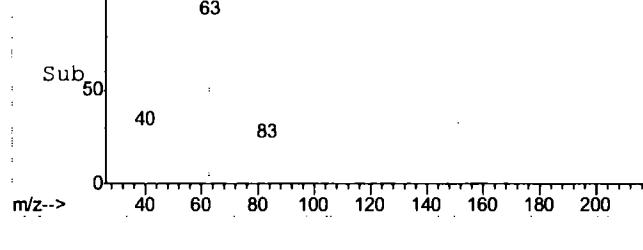
Tgt Ion: 58 Resp: 2209
Ion Ratio Lower Upper
58 100
43 304.1 268.4 328.4



#35
1,1-dichloroethane
Concen: 0.25 ug/L
RT: 7.45 min Scan# 778
Delta R.T. -0.01 min
Lab File: 4B54428.D
Acq: 12 Aug 2015 8:10 pm



Tgt Ion: 63 Resp: 1530
Ion Ratio Lower Upper
63 100
65 58.5 1.3 61.3
83 22.1 0.0 42.6



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54429.D
 Acq On : 12 Aug 2015 8:38 pm
 Operator : TOANP
 Sample : jc1107-6
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 25 Sample Multiplier: 1

Quant Time: Aug 13 11:56:56 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	145128	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	431535	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	484038	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	425202	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	215693	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	153306	49.13	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	98.26%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	161032	46.96	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	93.92%	
80) toluene-d8 (s)	11.02	98	565731	49.65	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	99.30%	
105) 4-bromofluorobenzene (s)	13.86	95	193473	51.38	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	102.76%	

Target Compounds

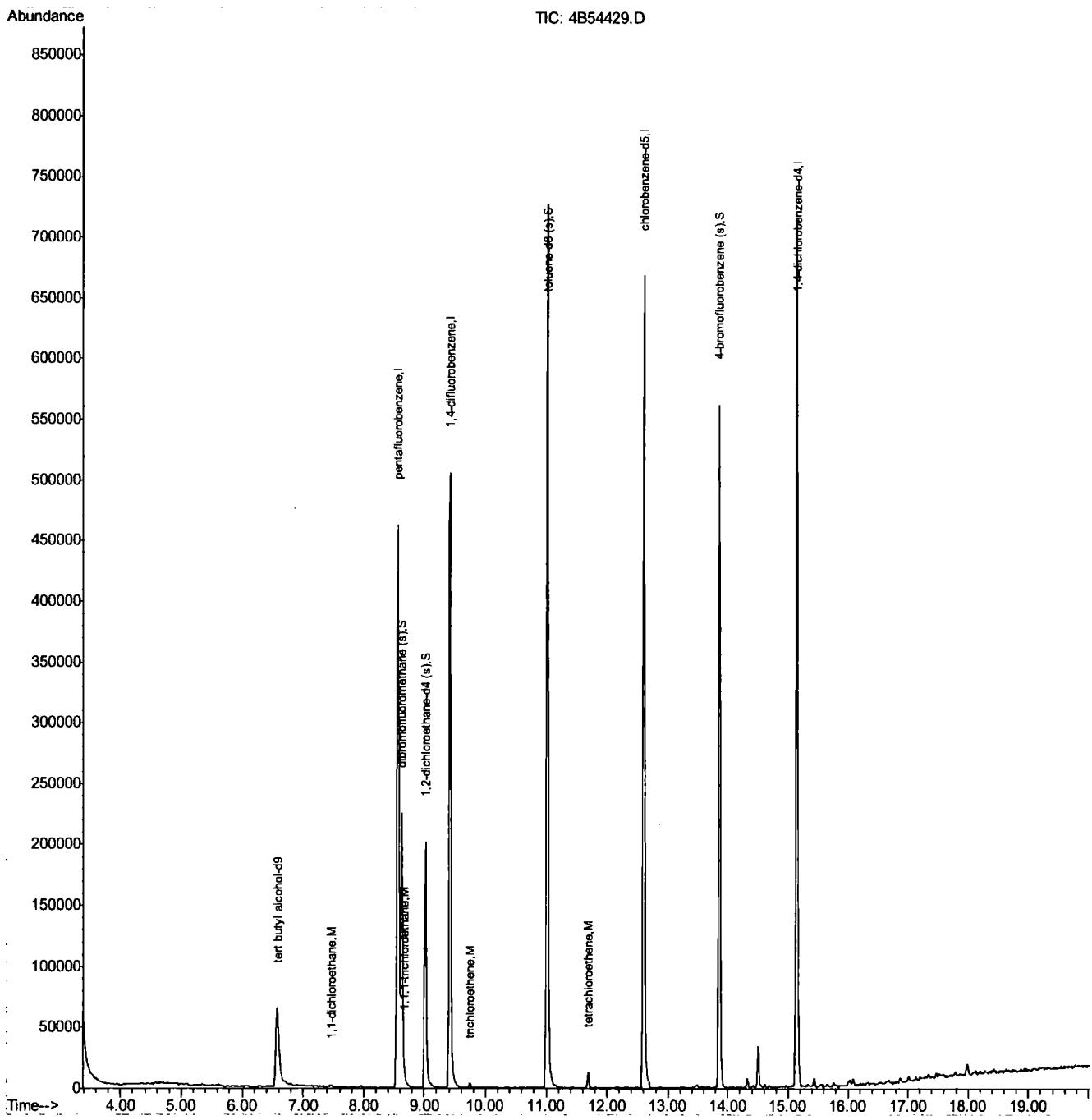
					Qvalue
35) 1,1-dichloroethane	7.45	63	2594	0.41	ug/L 75
53) 1,1,1-trichloroethane	8.67	97	1638	0.41	ug/L 78
69) trichloroethene	9.73	95	1495	0.44	ug/L 82
89) tetrachloroethene	11.69	164	3931	1.06	ug/L 90

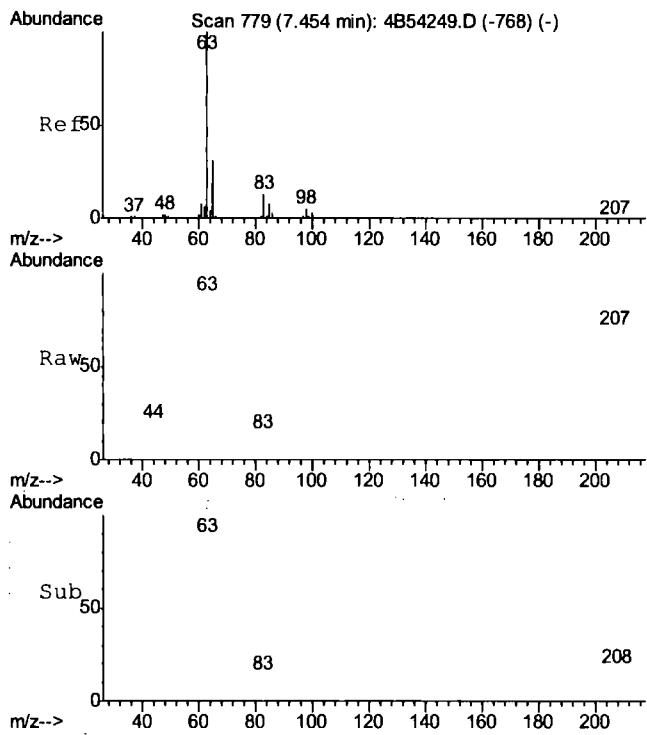
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54429.D
 Acq On : 12 Aug 2015 8:38 pm
 Operator : TOANP
 Sample : jc1107-6
 Misc : MS89470, V4B2296, w, , , 1
 ALS Vial : 25 Sample Multiplier: 1

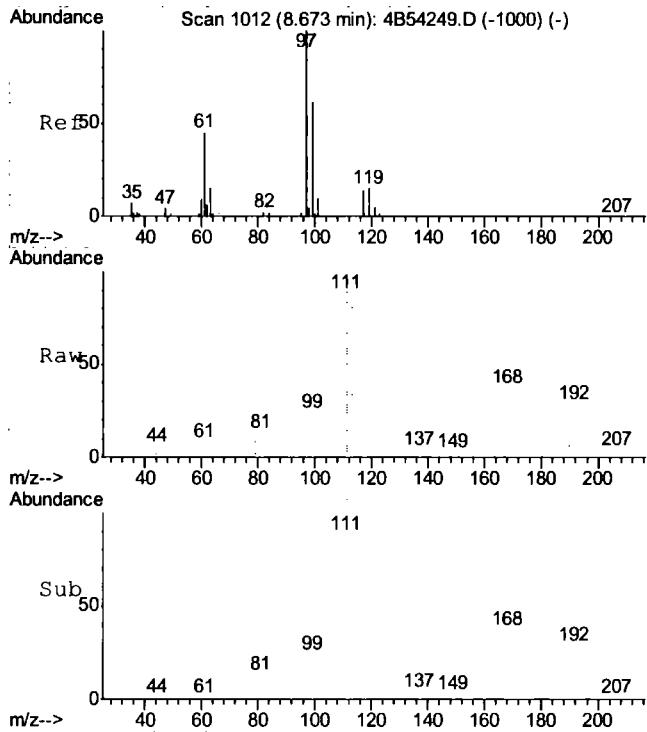
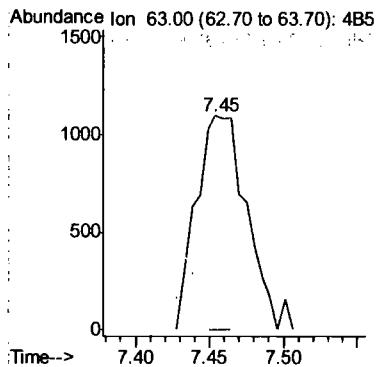
Quant Time: Aug 13 11:56:56 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration





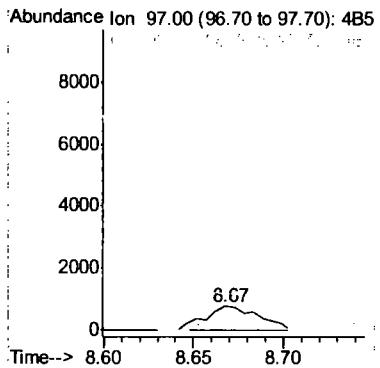
#35
1,1-dichloroethane
Concen: 0.41 ug/L
RT: 7.45 min Scan# 779
Delta R.T. -0.00 min
Lab File: 4B54429.D
Acq: 12 Aug 2015 8:38 pm

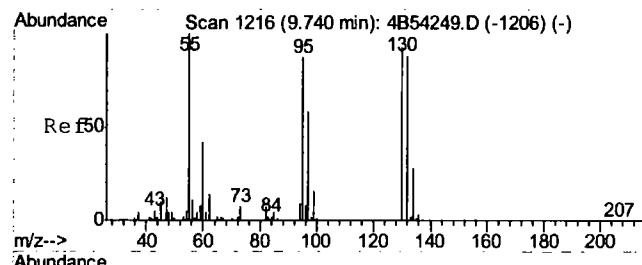
Tgt Ion: 63 Resp: 2594
Ion Ratio Lower Upper
63 100
65 50.2 1.3 61.3
83 14.0 0.0 42.6



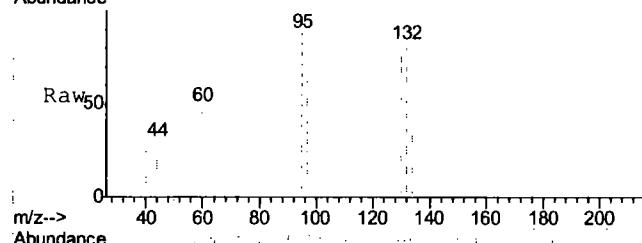
#53
1,1,1-trichloroethane
Concen: 0.41 ug/L
RT: 8.67 min Scan# 1011
Delta R.T. -0.01 min
Lab File: 4B54429.D
Acq: 12 Aug 2015 8:38 pm

Tgt Ion: 97 Resp: 1638
Ion Ratio Lower Upper
97 100
99 90.7 35.6 95.6
61 49.1 14.6 74.6





#69
trichloroethene
Concen: 0.44 ug/L
RT: 9.73 min Scan# 1215
Delta R.T. -0.01 min
Lab File: 4B54429.D
Acq: 12 Aug 2015 8:38 pm

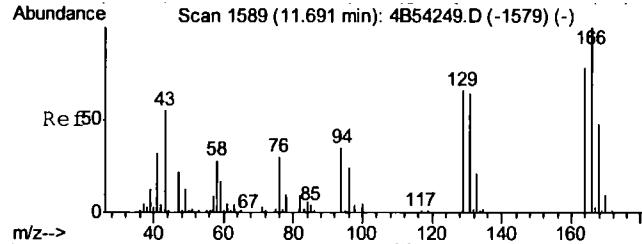
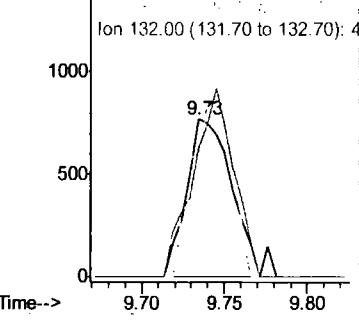
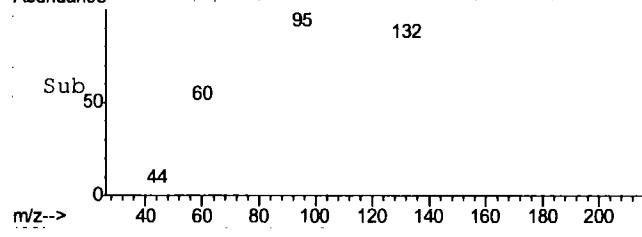


Tgt Ion: 95 Resp: 1495
Ion Ratio Lower Upper

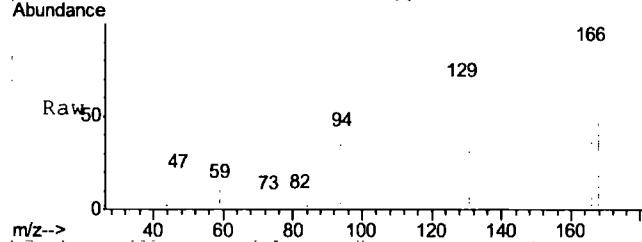
	95	100
97	67.1	36.5
130	78.4	76.6
132	81.5	71.3

Abundance Ion 95.00 (94.70 to 95.70): 4B5

Ion 132.00 (131.70 to 132.70): 4



#89
tetrachloroethene
Concen: 1.06 ug/L
RT: 11.69 min Scan# 1589
Delta R.T. -0.00 min
Lab File: 4B54429.D
Acq: 12 Aug 2015 8:38 pm

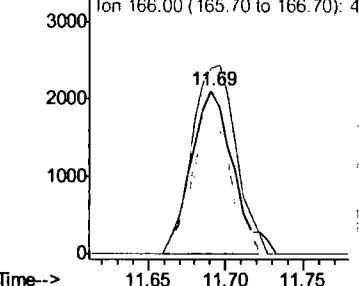
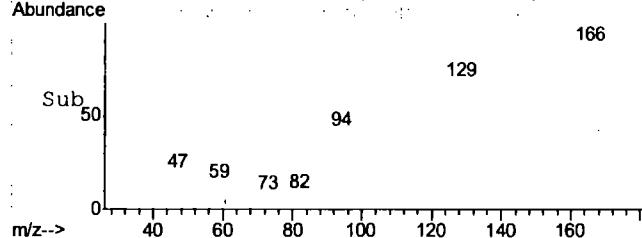


Tgt Ion: 164 Resp: 3931
Ion Ratio Lower Upper

	164	100
129	78.9	54.7
131	74.3	52.0
166	113.7	98.8

Abundance Ion 164.00 (163.70 to 164.70): 4

Ion 166.00 (165.70 to 166.70): 4



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54451.D
 Acq On : 13 Aug 2015 6:51 am
 Operator : TOANP
 Sample : jc1107-7
 Misc : MS89470,V4B2297,w,,,1
 ALS Vial : 47 Sample Multiplier: 1

Quant Time: Aug 13 12:27:21 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	181708	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	443640	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	499255	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	435289	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	212248	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	155837	48.58	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	97.16%
50) 1,2-dichloroethane-d4 (s)	9.02	65	161936	45.93	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	91.86%
80) toluene-d8 (s)	11.02	98	576442	49.05	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	98.10%
105) 4-bromofluorobenzene (s)	13.86	95	193363	52.18	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.36%

Target Compounds

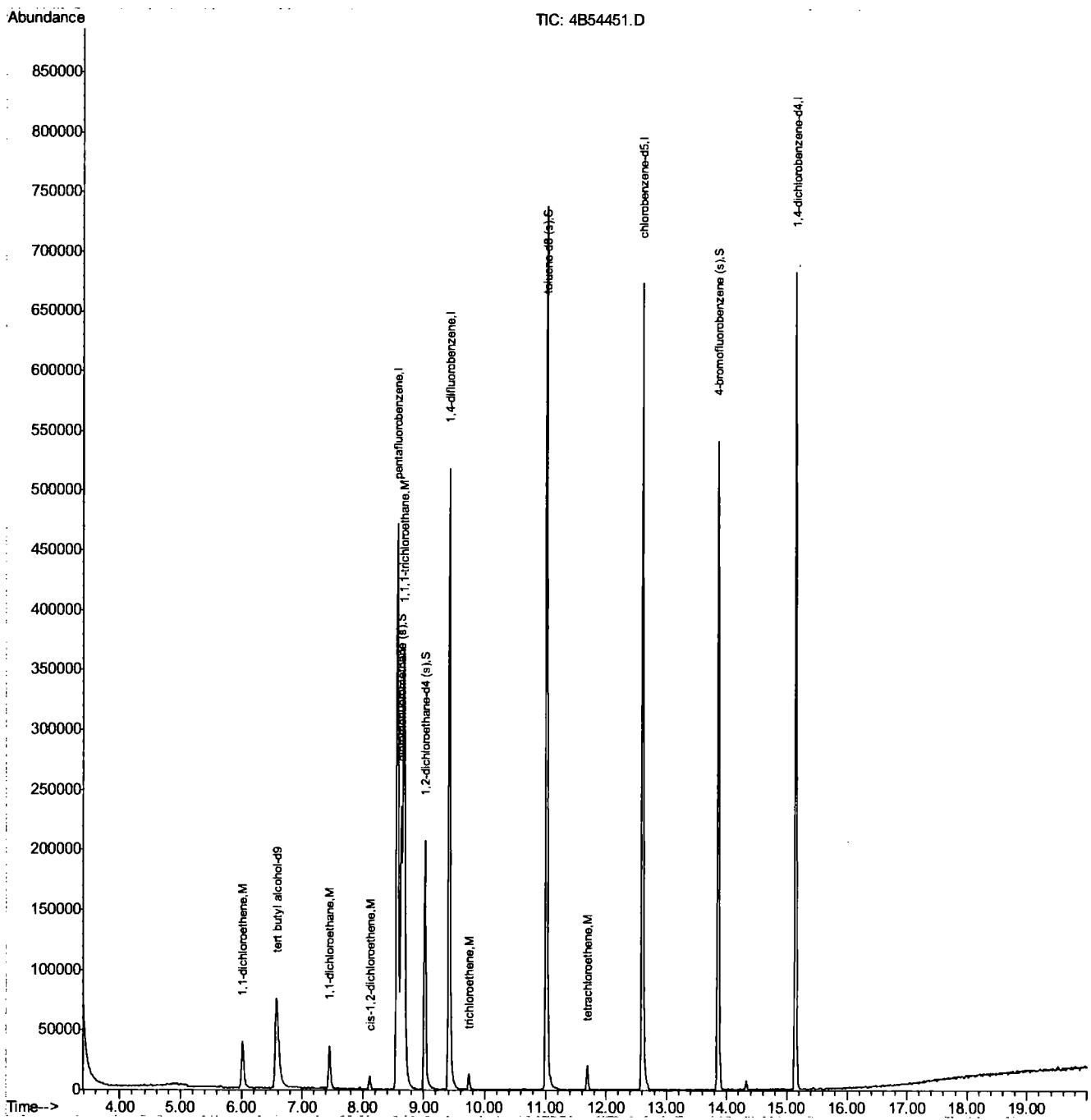
					Qvalue
22) 1,1-dichloroethene	6.02	96	24240	6.92	ug/L
35) 1,1-dichloroethane	7.46	63	46438	7.17	ug/L
42) cis-1,2-dichloroethene	8.11	96	5968	1.59	ug/L
53) 1,1,1-trichloroethane	8.67	97	311978	76.16	ug/L
69) trichloroethene	9.74	95	5132	1.46	ug/L
89) tetrachloroethene	11.69	164	5791	1.53	ug/L

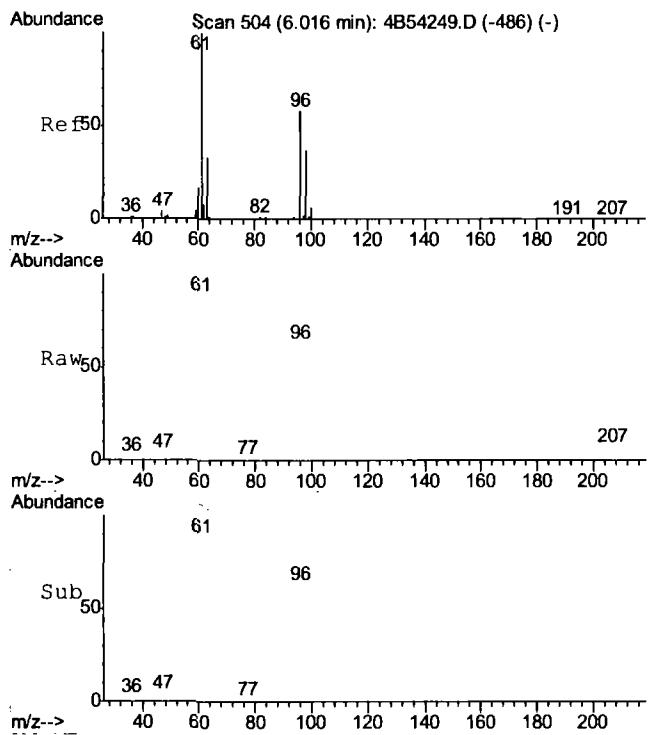
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54451.D
 Acq On : 13 Aug 2015 6:51 am
 Operator : TOANP
 Sample : jc1107-7
 Misc : MS89470, V4B2297, w,,,1
 ALS Vial : 47 Sample Multiplier: 1

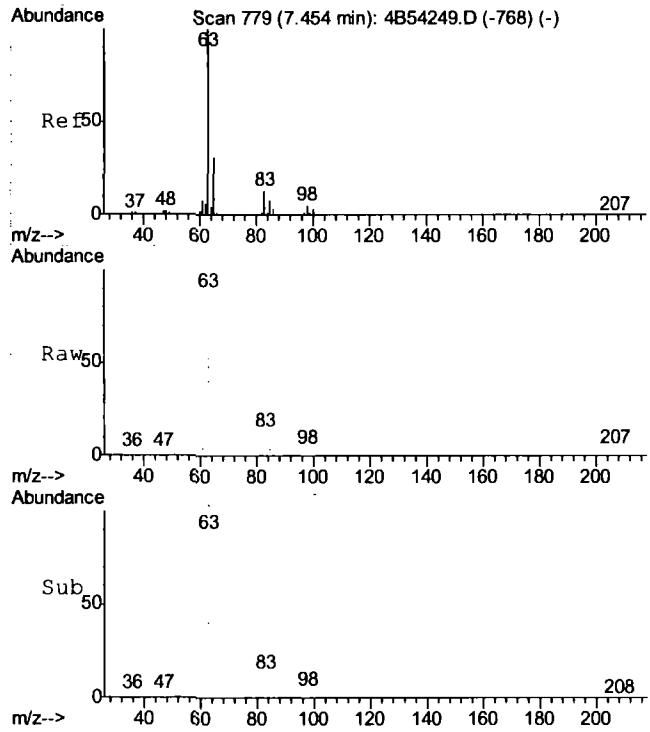
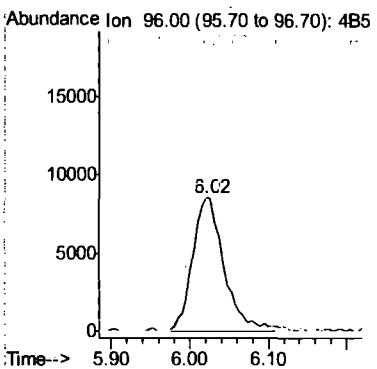
Quant Time: Aug 13 12:27:21 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration





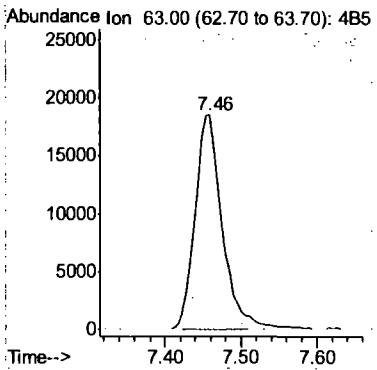
#22
1,1-dichloroethene
Concen: 6.92 ug/L
RT: 6.02 min Scan# 505
Delta R.T. 0.01 min
Lab File: 4B54451.D
Acq: 13 Aug 2015 6:51 am

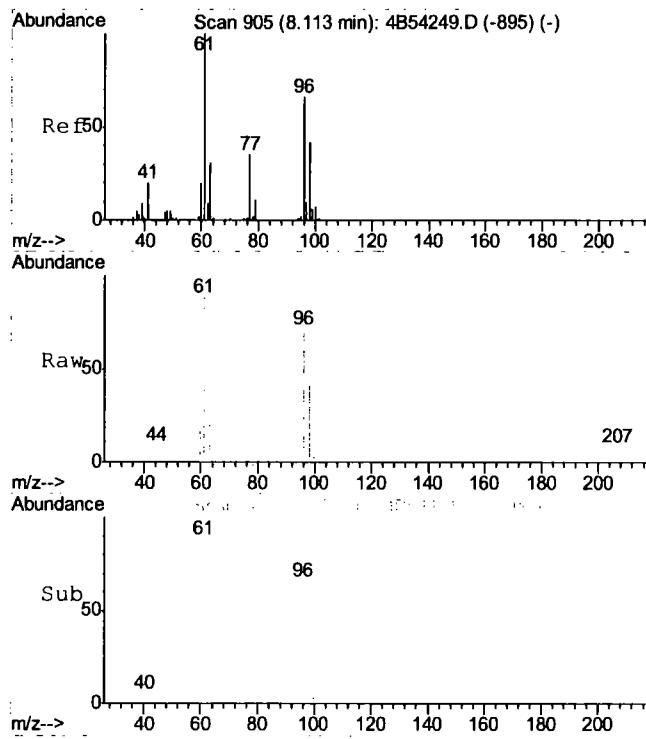
Tgt Ion: 96 Resp: 24240
Ion Ratio Lower Upper
96 100
61 159.7 141.9 201.9
63 50.8 26.8 86.8



#35
1,1-dichloroethane
Concen: 7.17 ug/L
RT: 7.46 min Scan# 780
Delta R.T. 0.01 min
Lab File: 4B54451.D
Acq: 13 Aug 2015 6:51 am

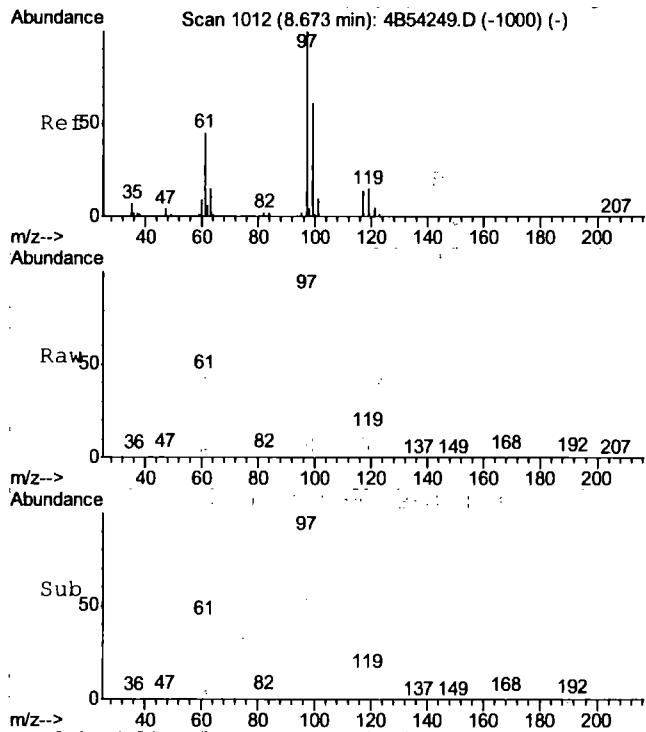
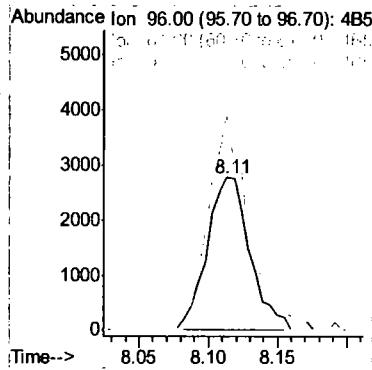
Tgt Ion: 63 Resp: 46438
Ion Ratio Lower Upper
63 100
65 33.9 1.3 61.3
83 13.0 0.0 42.6





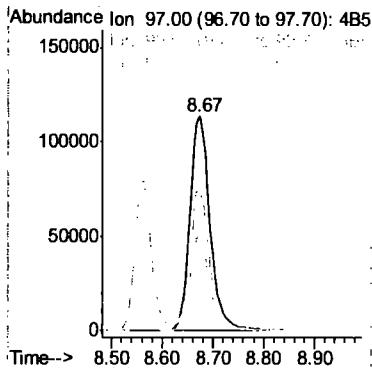
#42
cis-1,2-dichloroethene
Concen: 1.59 ug/L
RT: 8.11 min Scan# 905
Delta R.T. -0.00 min
Lab File: 4B54451.D
Acq: 13 Aug 2015 6:51 am

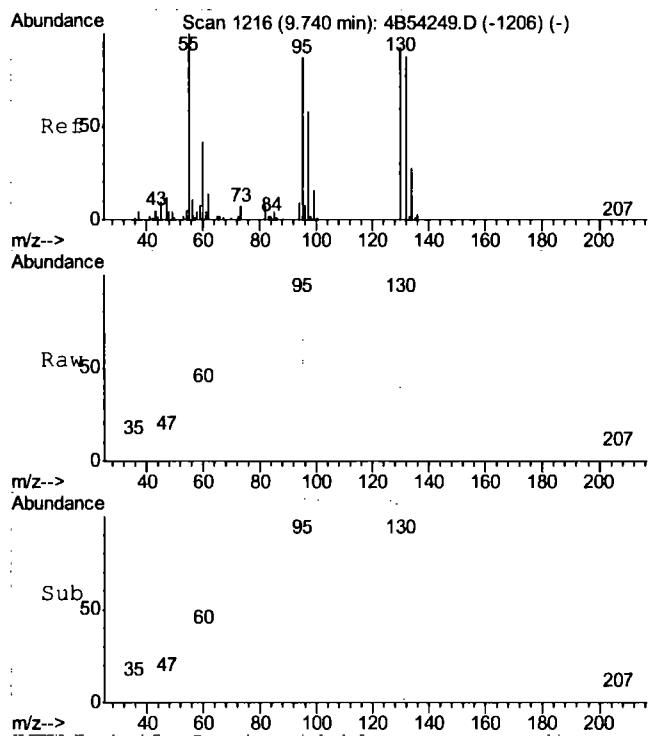
Tgt Ion: 96 Resp: 5968
Ion Ratio Lower Upper
96 100
61 140.6 123.9 183.9
98 60.5 34.0 94.0



#53
1,1,1-trichloroethane
Concen: 76.16 ug/L
RT: 8.67 min Scan# 1012
Delta R.T. -0.00 min
Lab File: 4B54451.D
Acq: 13 Aug 2015 6:51 am

Tgt Ion: 97 Resp: 311978
Ion Ratio Lower Upper
97 100
99 66.6 35.6 95.6
61 45.4 14.6 74.6

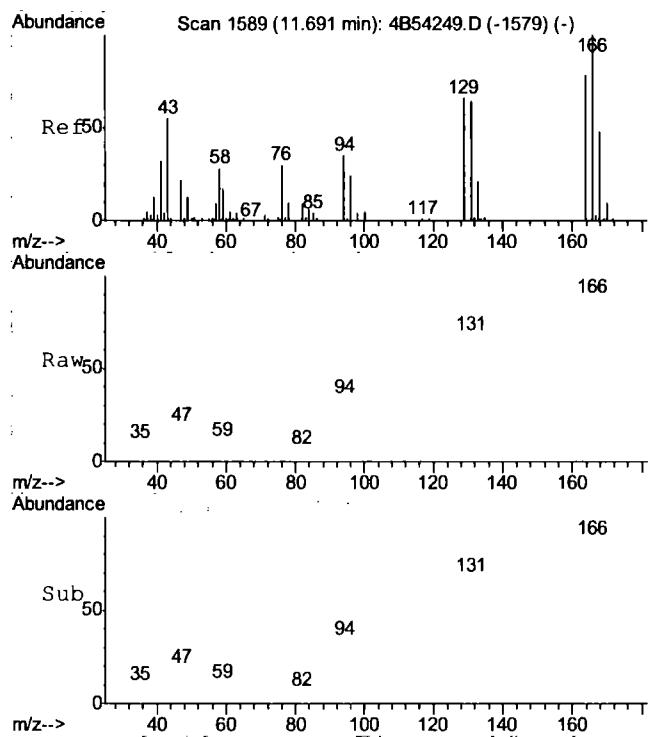
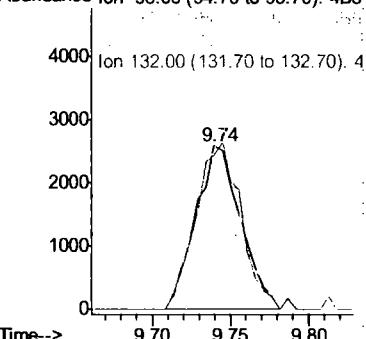




#69
trichloroethene
Concen: 1.46 ug/L
RT: 9.74 min Scan# 1216
Delta R.T. -0.00 min
Lab File: 4B54451.D
Acq: 13 Aug 2015 6:51 am

Tgt Ion: 95 Resp: 5132
Ion Ratio Lower Upper
95 100
97 74.1 36.5 96.5
130 96.7 76.6 136.6
132 94.1 71.3 131.3

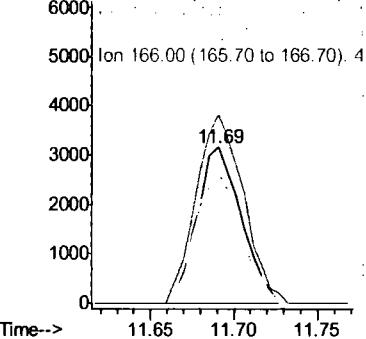
Abundance Ion 95.00 (94.70 to 95.70): 4B5



#89
tetrachloroethene
Concen: 1.53 ug/L
RT: 11.69 min Scan# 1589
Delta R.T. -0.00 min
Lab File: 4B54451.D
Acq: 13 Aug 2015 6:51 am

Tgt Ion: 164 Resp: 5791
Ion Ratio Lower Upper
164 100
129 76.6 54.7 114.7
131 82.3 52.0 112.0
166 121.3 98.8 158.8

Abundance Ion 164.00 (163.70 to 164.70): 4



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54450.D
 Acq On : 13 Aug 2015 6:23 am
 Operator : TOANP
 Sample : jc1107-8
 Misc : MS89470,V4B2297,w,,,1
 ALS Vial : 46 Sample Multiplier: 1

Quant Time: Aug 13 12:26:41 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	144786	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	464734	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	523670	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	454873	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	216894	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	163277	48.59	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery =	97.18%		
50) 1,2-dichloroethane-d4 (s)	9.02	65	168285	45.57	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery =	91.14%		
80) toluene-d8 (s)	11.02	98	606079	49.16	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery =	98.32%		
105) 4-bromofluorobenzene (s)	13.86	95	202524	53.48	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	106.96%		

Target Compounds

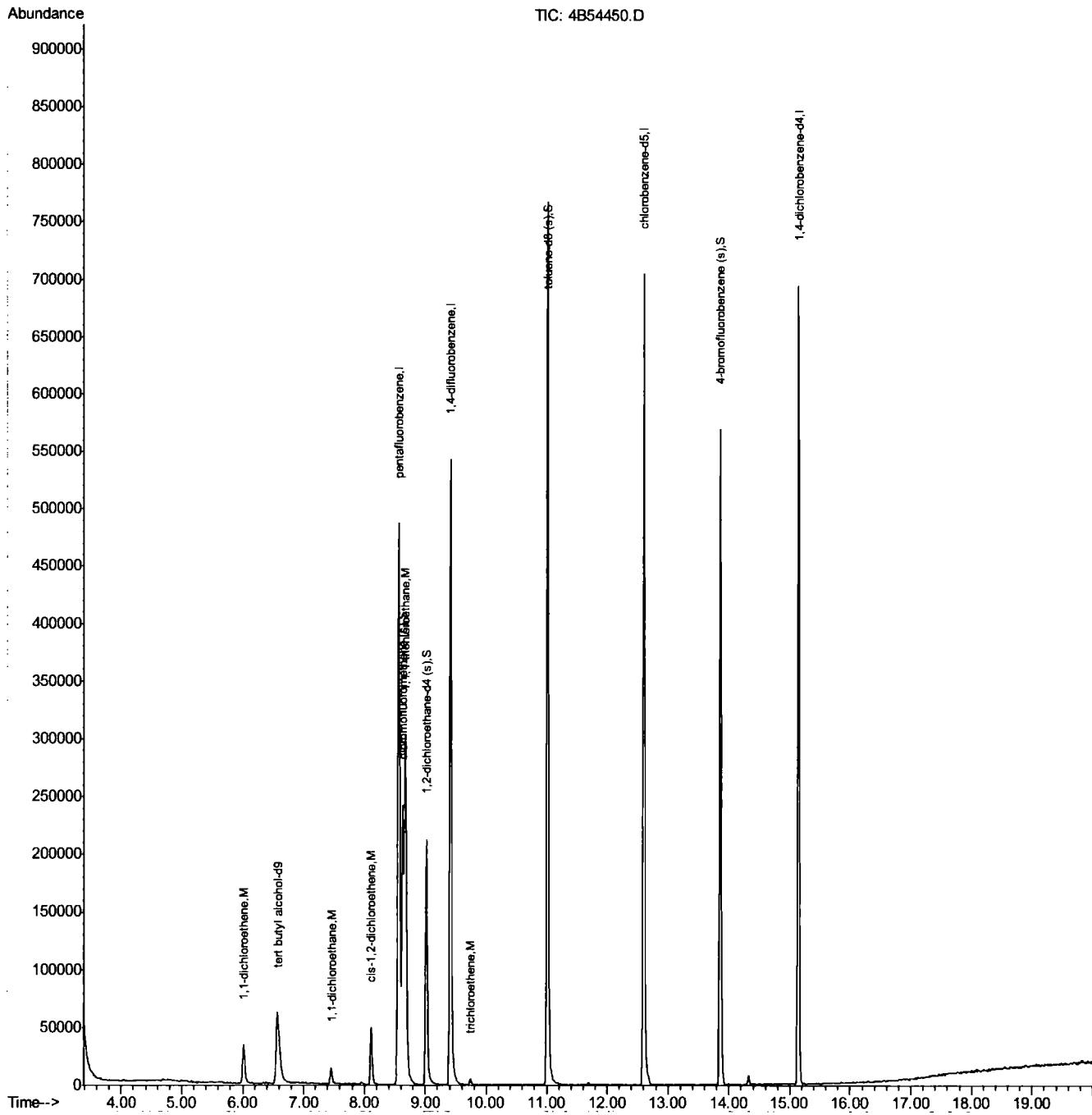
				Qvalue
22) 1,1-dichloroethene	6.03	96	20483	5.58 ug/L 89
35) 1,1-dichloroethane	7.46	63	17853	2.63 ug/L 97
42) cis-1,2-dichloroethene	8.11	96	26728	6.82 ug/L 90
53) 1,1,1-trichloroethane	8.67	97	255683	59.59 ug/L 97
69) trichloroethene	9.74	95	2248	0.61 ug/L 87

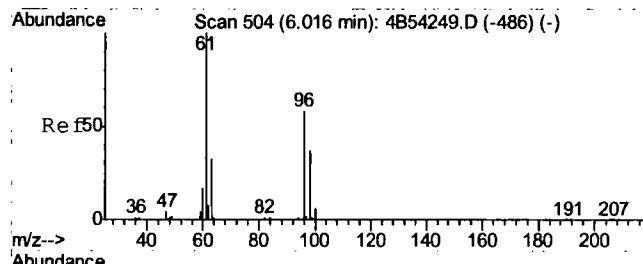
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

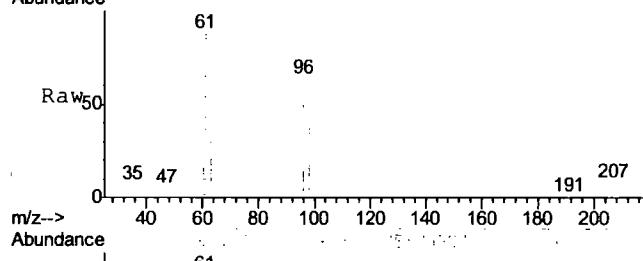
Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54450.D
 Acq On : 13 Aug 2015 6:23 am
 Operator : TOANP
 Sample : jc1107-8
 Misc : MS89470,V4B2297,w,,,1
 ALS Vial : 46 Sample Multiplier: 1

Quant Time: Aug 13 12:26:41 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

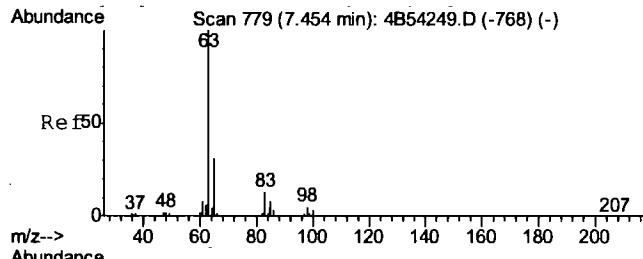
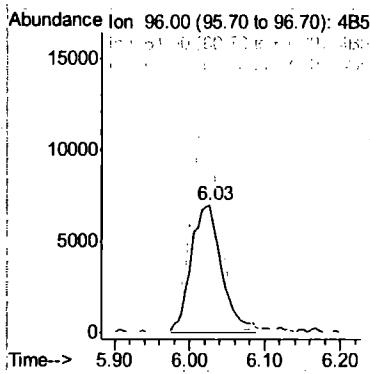
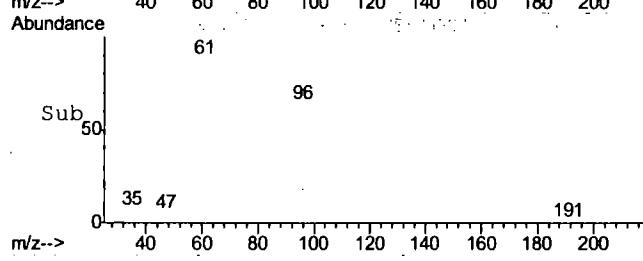




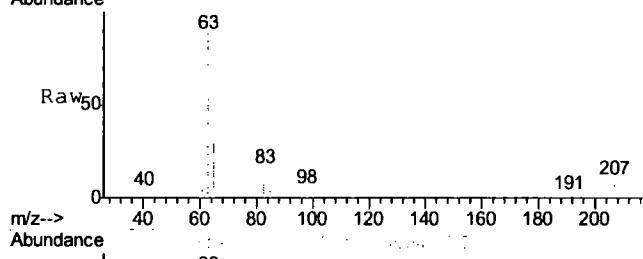
#22
1,1-dichloroethene
Concen: 5.58 ug/L
RT: 6.03 min Scan# 506
Delta R.T. 0.01 min
Lab File: 4B54450.D
Acq: 13 Aug 2015 6:23 am



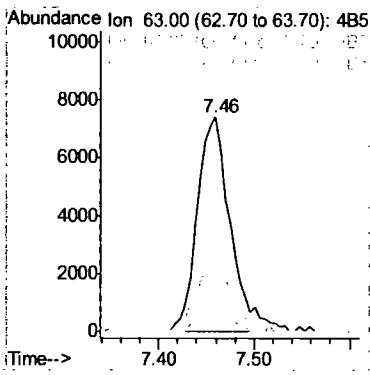
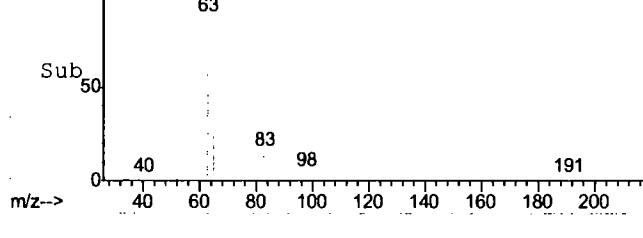
Tgt Ion: 96 Resp: 20483
Ion Ratio Lower Upper
96 100
61 156.9 141.9 201.9
63 48.8 26.8 86.8

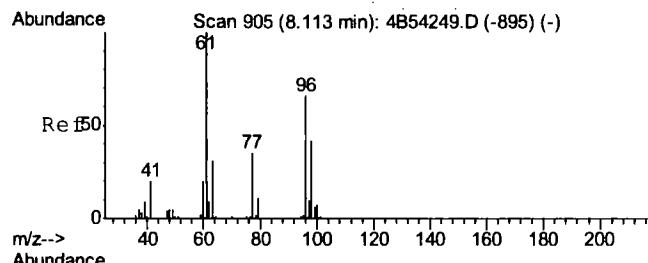


#35
1,1-dichloroethane
Concen: 2.63 ug/L
RT: 7.46 min Scan# 780
Delta R.T. 0.01 min
Lab File: 4B54450.D
Acq: 13 Aug 2015 6:23 am

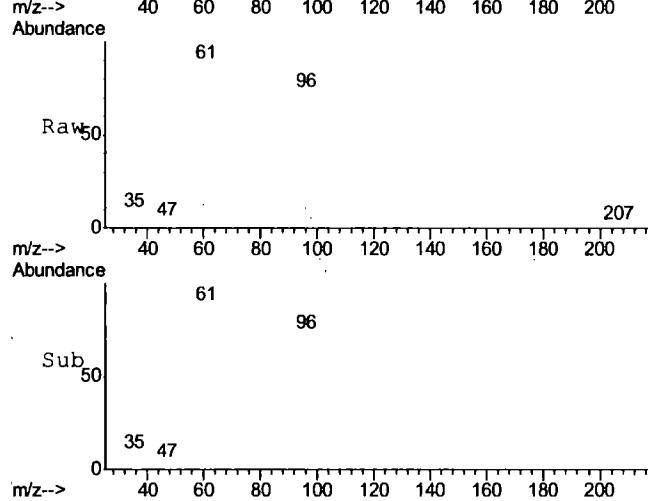


Tgt Ion: 63 Resp: 17853
Ion Ratio Lower Upper
63 100
65 30.9 1.3 61.3
83 15.5 0.0 42.6

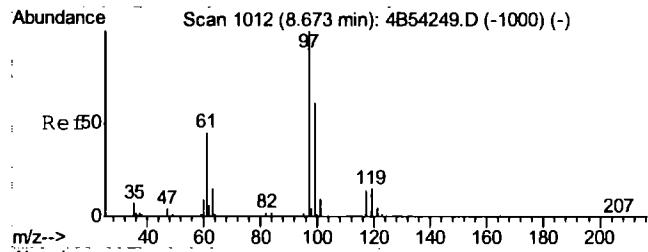




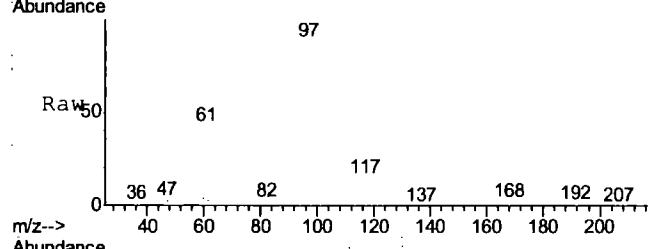
#42
cis-1,2-dichloroethene
Concen: 6.82 ug/L
RT: 8.11 min Scan# 904
Delta R.T. -0.01 min
Lab File: 4B54450.D
Acq: 13 Aug 2015 6:23 am



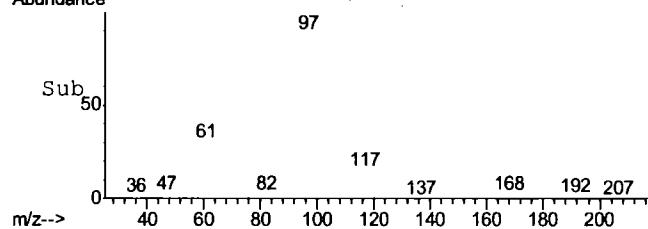
Tgt Ion: 96 Resp: 26728
Ion Ratio Lower Upper
96 100
61 136.8 123.9 183.9
98 61.6 34.0 94.0



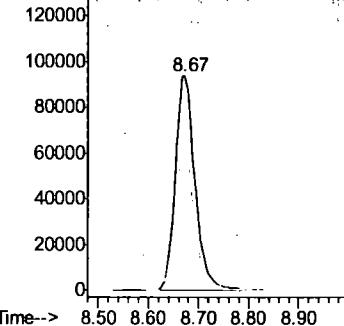
#53
1,1,1-trichloroethane
Concen: 59.59 ug/L
RT: 8.67 min Scan# 1012
Delta R.T. -0.00 min
Lab File: 4B54450.D
Acq: 13 Aug 2015 6:23 am

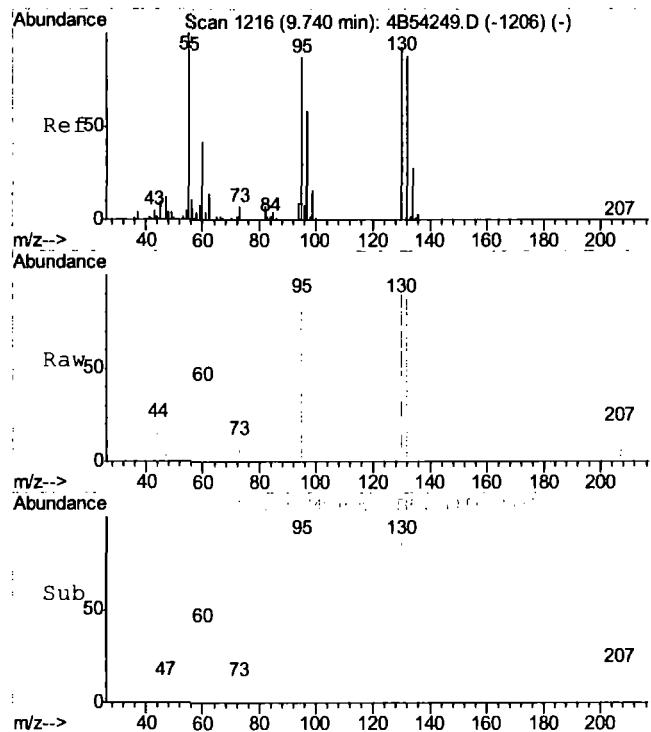


Tgt Ion: 97 Resp: 255683
Ion Ratio Lower Upper
97 100
99 67.9 35.6 95.6
61 42.8 14.6 74.6



Abundance Ion 97.00 (96.70 to 97.70): 4B5





69
trichloroethene
Concen: 0.61 ug/L
RT: 9.74 min Scan# 1216
Delta R.T. -0.00 min
Lab File: 4B54450.D
Acq: 13 Aug 2015 6:23 am

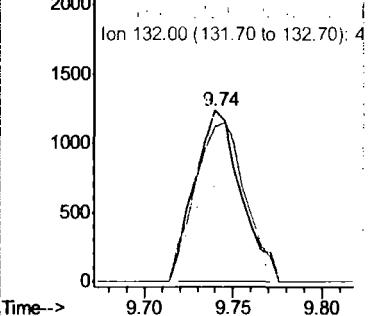
Tgt Ion: 95 Resp: 2248

Ion Ratio Lower Upper

95	100		
97	56.1	36.5	96.5
130	91.2	76.6	136.6
132	90.1	71.3	131.3

Abundance Ion 95.00 (94.70 to 95.70): 4B5

Ion 132.00 (131.70 to 132.70): 4B5



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54452.D
 Acq On : 13 Aug 2015 7:19 am
 Operator : TOANP
 Sample : jc1107-9
 Misc : MS89470,V4B2297,w,,,1
 ALS Vial : 48 Sample Multiplier: 1

Quant Time: Aug 13 12:27:52 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	142504	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	438675	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	494969	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	432890	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	208348	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	153512	48.40	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery =	96.80%		
50) 1,2-dichloroethane-d4 (s)	9.02	65	162448	46.60	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery =	93.20%		
80) toluene-d8 (s)	11.02	98	573762	49.24	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery =	98.48%		
105) 4-bromofluorobenzene (s)	13.86	95	192346	52.88	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	105.76%		

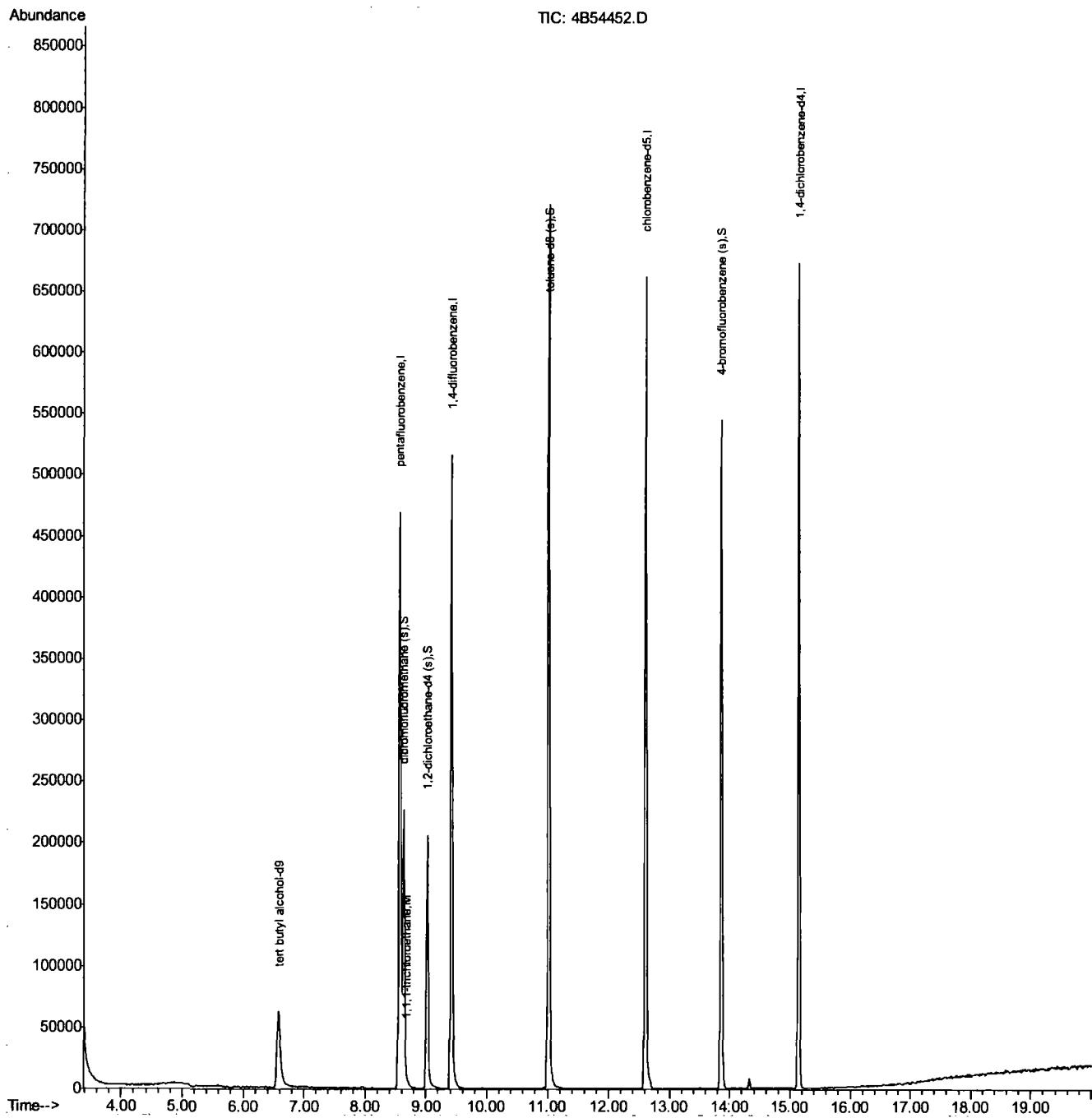
Target Compounds				Qvalue	
53) 1,1,1-trichloroethane	8.67	97	2089	0.52	ug/L # 63

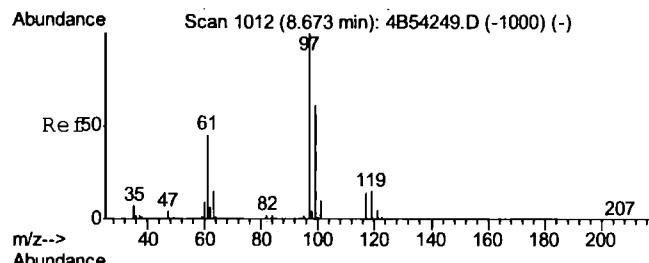
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54452.D
Acq On : 13 Aug 2015 7:19 am
Operator : TOANP
Sample : jc1107-9
Misc : MS89470, V4B2297, w, , , , 1
ALS Vial : 48 Sample Multiplier: 1

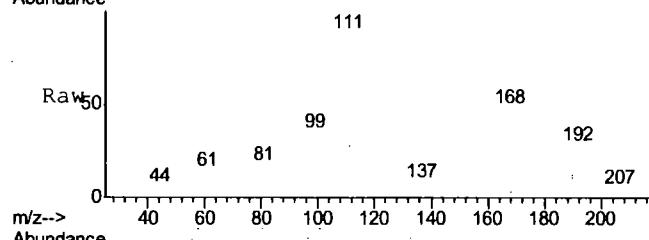
Quant Time: Aug 13 12:27:52 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



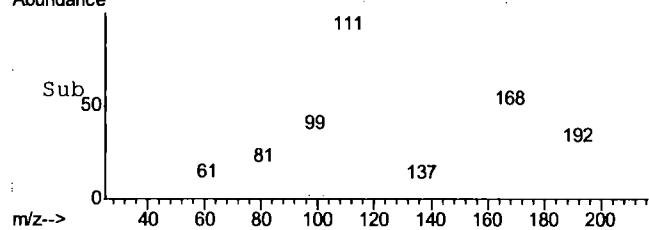
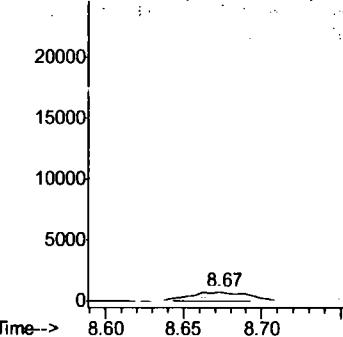


#53
1,1,1-trichloroethane
Concen: 0.52 ug/L
RT: 8.67 min Scan# 1012
Delta R.T. -0.00 min
Lab File: 4B54452.D
Acq: 13 Aug 2015 7:19 am

Tgt Ion: 97 Resp: 2089
Ion Ratio Lower Upper
97 100
99 100.7 35.6 95.6#
61 60.9 14.6 74.6



Abundance Ion 97.00 (96.70 to 97.70): 4B5



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54453.D
 Acq On : 13 Aug 2015 7:48 am
 Operator : TOANP
 Sample : jc1107-10
 Misc : MS89470,V4B2297,w,,,,,1
 ALS Vial : 49 Sample Multiplier: 1

Quant Time: Aug 13 12:28:23 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	134311	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	433187	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	481637	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	421297	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	205804	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	152759	48.77	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 97.54%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	161927	47.04	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 94.08%	
80) toluene-d8 (s)	11.02	98	565751	49.90	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 99.80%	
105) 4-bromofluorobenzene (s)	13.86	95	188288	52.40	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 104.80%	

Target Compounds

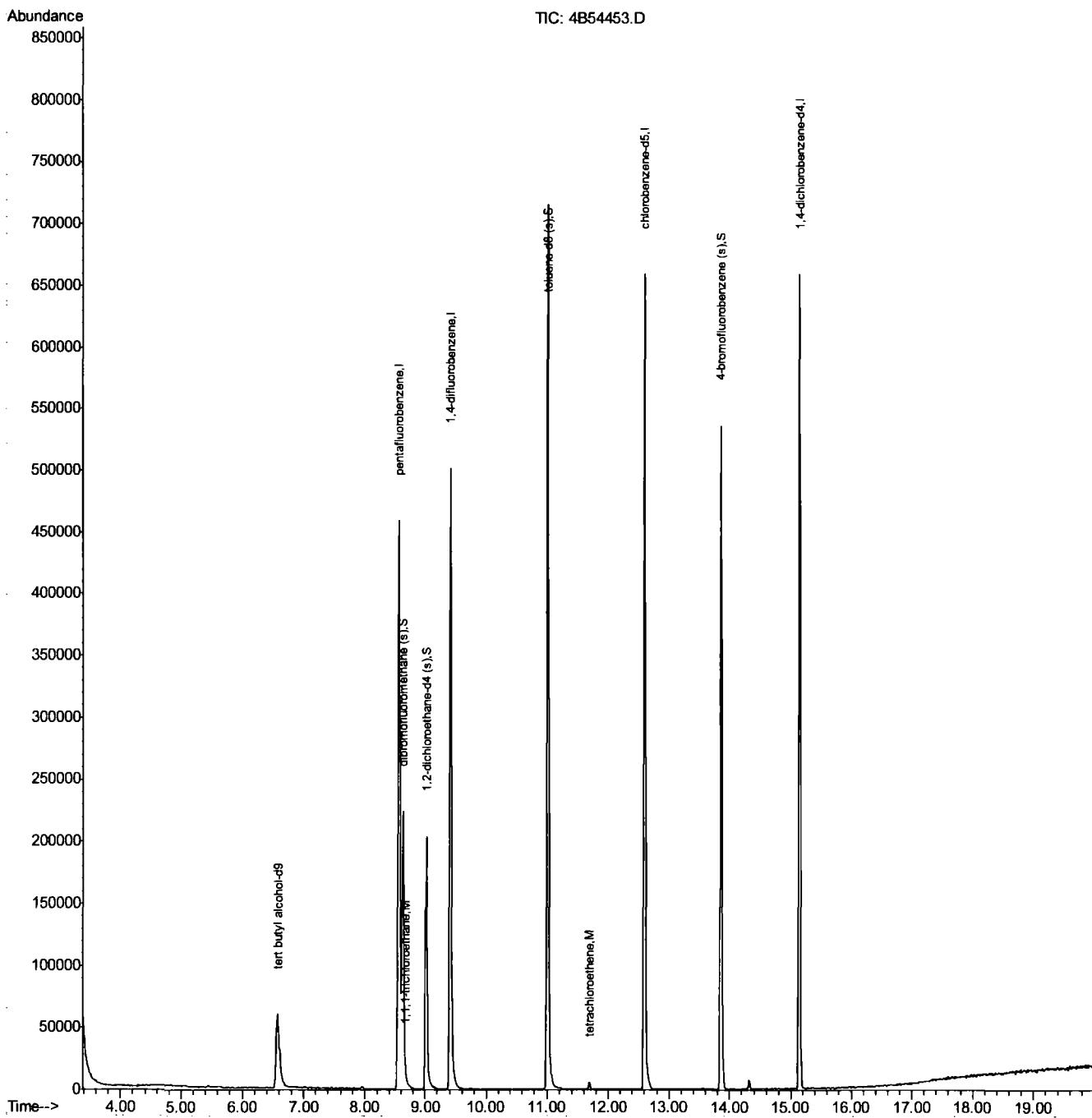
				Qvalue
53) 1,1,1-trichloroethane	8.68	97	1563	0.39 ug/L # 64
89) tetrachloroethene	11.69	164	1592	0.43 ug/L # 79

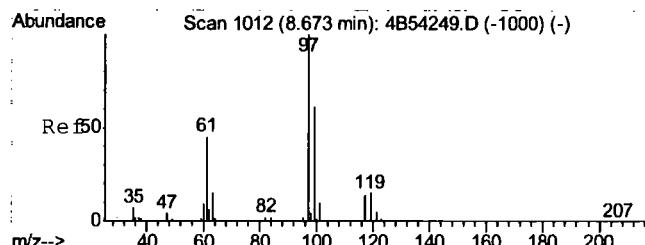
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

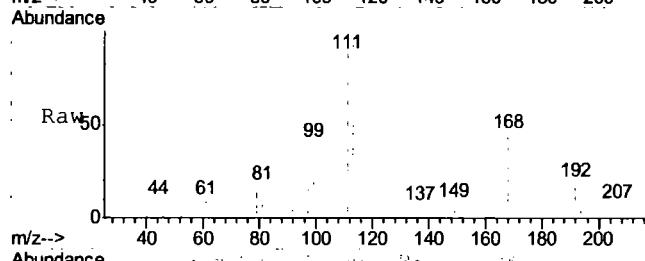
Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54453.D
Acq On : 13 Aug 2015 7:48 am
Operator : TOANP
Sample : jc1107-10
Misc : MS89470,V4B2297,w,,,1
ALS Vial : 49 Sample Multiplier: 1

Quant Time: Aug 13 12:28:23 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration

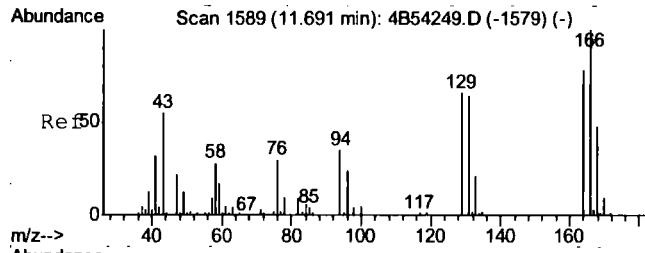
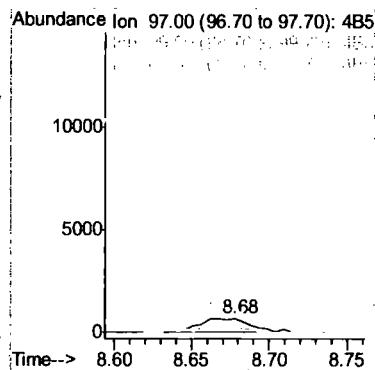
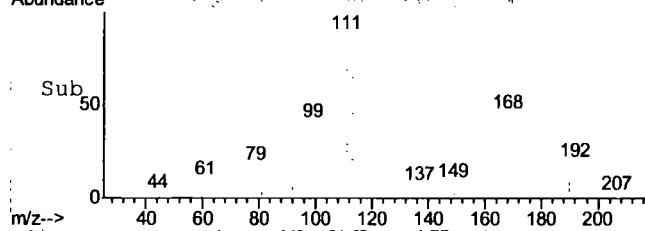




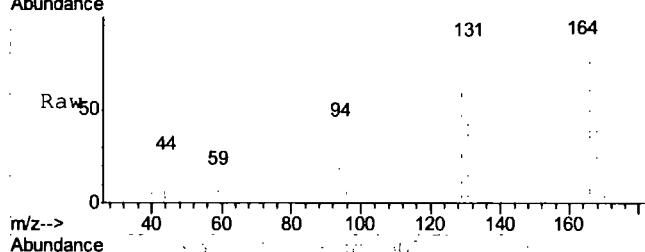
#53
1,1,1-trichloroethane
Concen: 0.39 ug/L
RT: 8.68 min Scan# 1013
Delta R.T. 0.00 min
Lab File: 4B54453.D
Acq: 13 Aug 2015 7:48 am



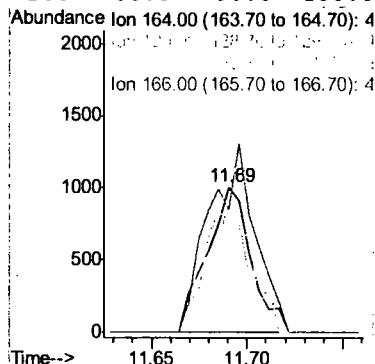
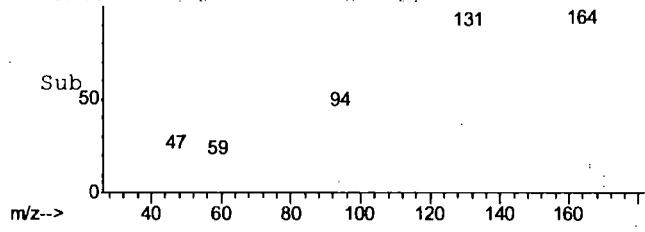
Tgt Ion: 97 Resp: 1563
Ion Ratio Lower Upper
97 100
99 109.5 35.6 95.6#
61 40.3 14.6 74.6



#89
tetrachloroethene
Concen: 0.43 ug/L
RT: 11.69 min Scan# 1589
Delta R.T. -0.00 min
Lab File: 4B54453.D
Acq: 13 Aug 2015 7:48 am



Tgt Ion: 164 Resp: 1592
Ion Ratio Lower Upper
164 100
129 77.1 54.7 114.7
131 87.2 52.0 112.0
166 83.8 98.8 158.8#



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54423.D
 Acq On : 12 Aug 2015 5:51 pm
 Operator : TOANP
 Sample : jc1107-11
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 19 Sample Multiplier: 1

Quant Time: Aug 13 11:51:56 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	140434	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	450428	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	498750	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	437102	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	216722	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	155872	47.86	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	95.72%
50) 1,2-dichloroethane-d4 (s)	9.02	65	163515	45.68	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	91.36%
80) toluene-d8 (s)	11.02	98	584968	49.82	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	99.64%
105) 4-bromofluorobenzene (s)	13.86	95	196797	52.01	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.02%

Target Compounds

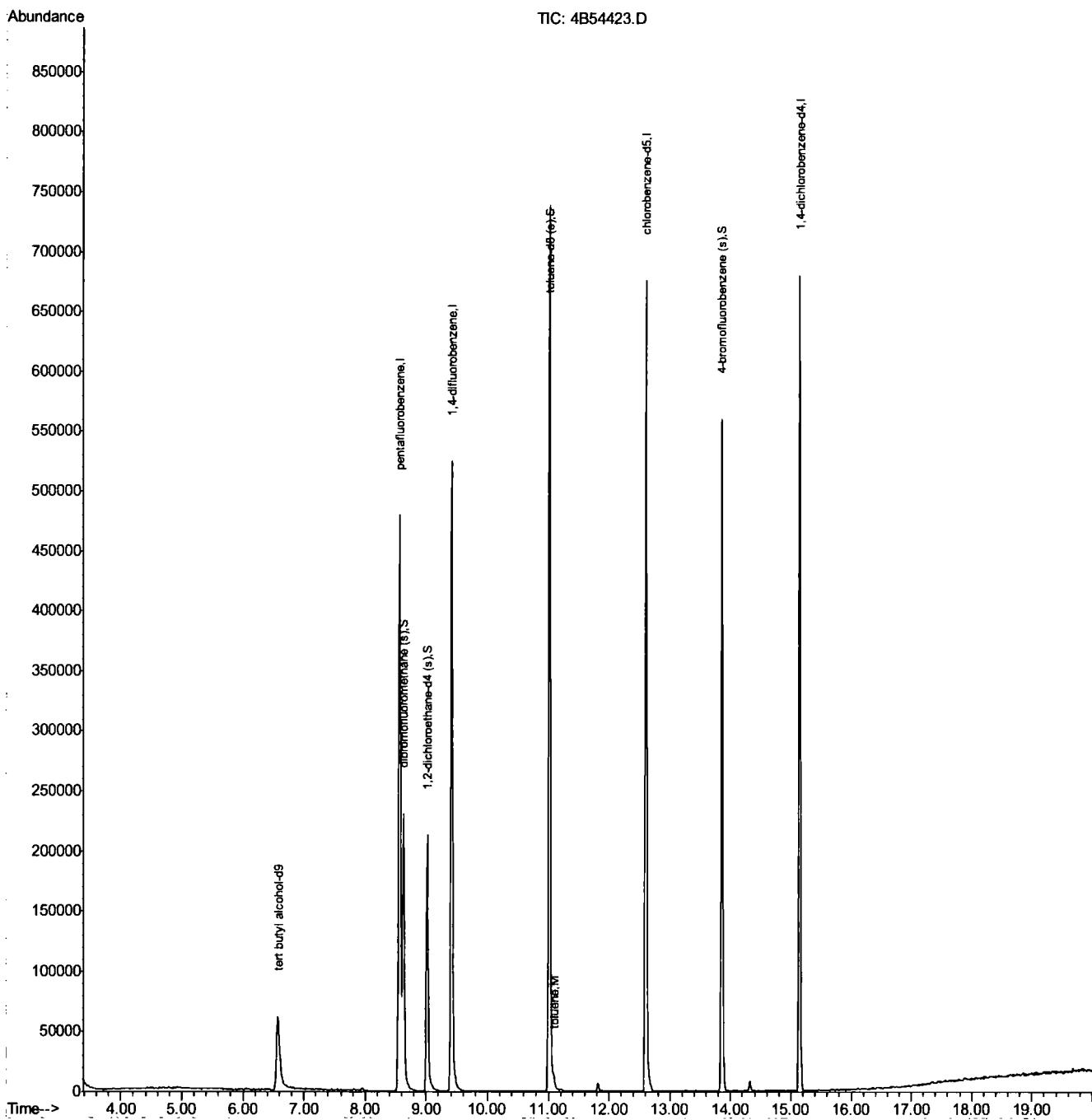
82) toluene	11.09	92	2632	0.32	ug/L	Qvalue 88
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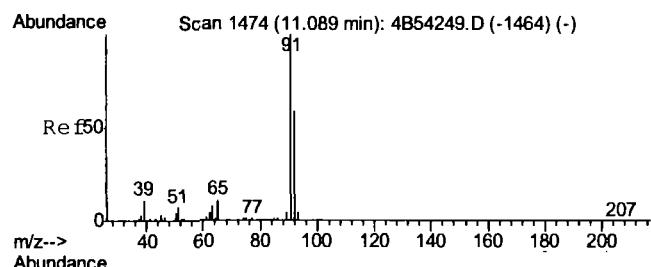
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54423.D
Acq On : 12 Aug 2015 5:51 pm
Operator : TOANP
Sample : jc1107-11
Misc : MS89470, V4B2296, w, , , 1
ALS Vial : 19 Sample Multiplier: 1

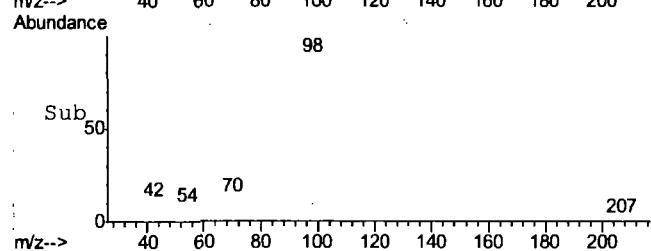
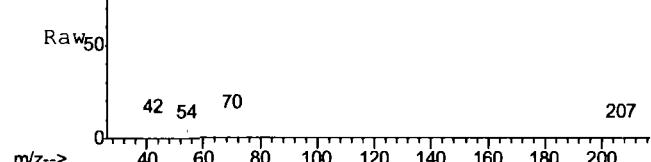
Quant Time: Aug 13 11:51:56 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



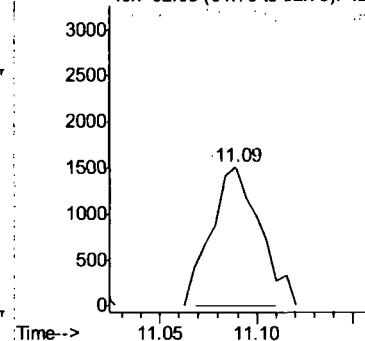


#82
toluene
Concen: 0.32 ug/L
RT: 11.09 min Scan# 1474
Delta R.T. -0.00 min
Lab File: 4B54423.D
Acq: 12 Aug 2015 5:51 pm

Tgt Ion: 92 Resp: 2632
Ion Ratio Lower Upper
92 100
91 153.7 140.0 200.0
65 26.4 0.0 49.3



Abundance Ion 92.00 (91.70 to 92.70): 4B5



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54424.D
 Acq On : 12 Aug 2015 6:18 pm
 Operator : TOANP
 Sample : jc1107-12
 Misc : MS89470,V4B2296,w,,,,,1
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 13 11:52:30 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	137119	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	445018	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	501240	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	439157	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	212599	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	154156	47.91	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	95.82%
50) 1,2-dichloroethane-d4 (s)	9.02	65	163125	46.13	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	92.26%
80) toluene-d8 (s)	11.02	98	583978	49.49	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	98.98%
105) 4-bromofluorobenzene (s)	13.86	95	195413	52.65	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	105.30%

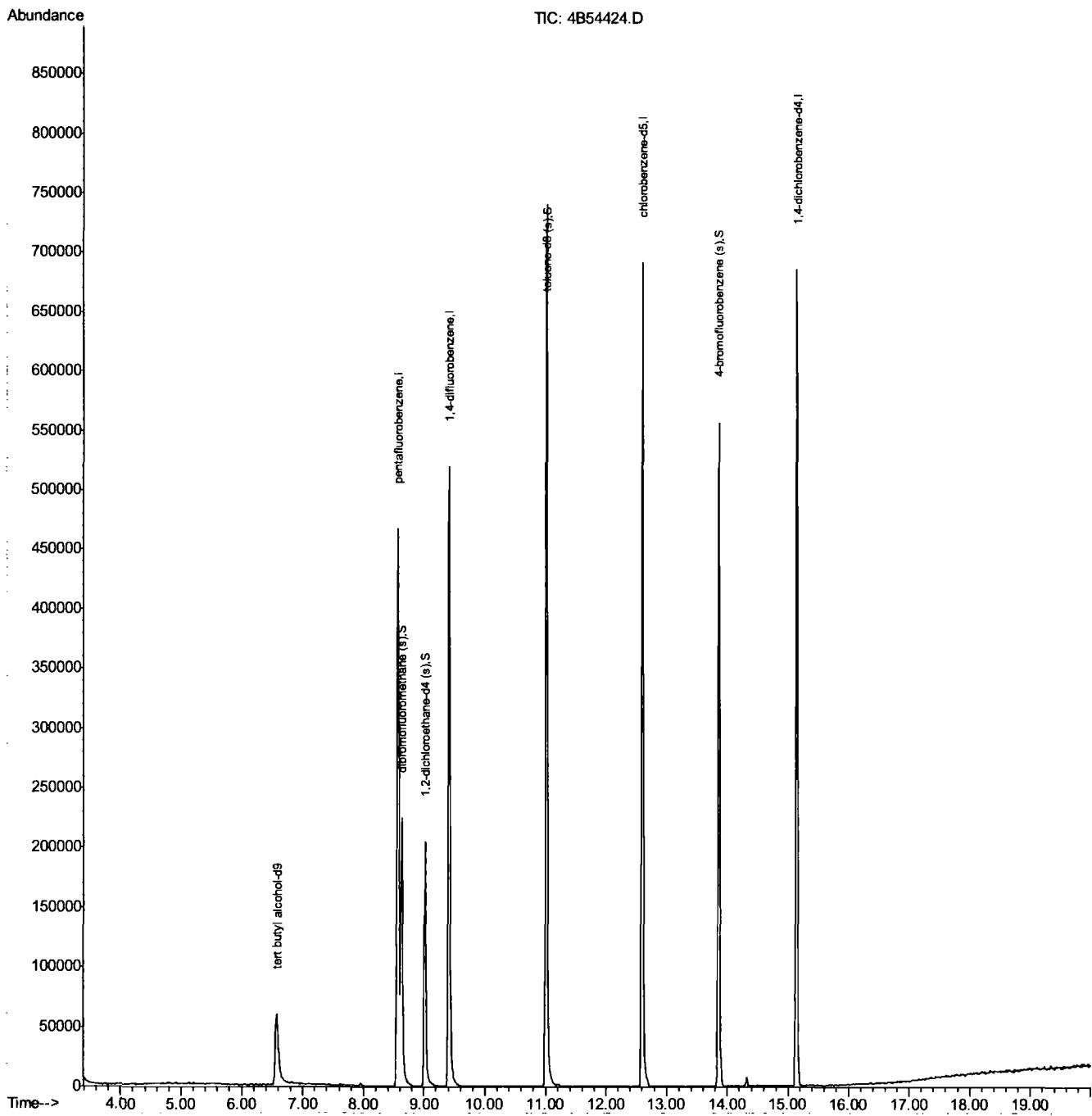
Target Compounds	Qvalue
(#)	= qualifier out of range (m) = manual integration (+) = signals summed

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54424.D
 Acq On : 12 Aug 2015 6:18 pm
 Operator : TOANP
 Sample : jc1107-12
 Misc : MS89470,V4B2296,w,,,,,1
 ALS Vial : 20 Sample Multiplier: 1

Quant Time: Aug 13 11:52:30 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54409.D
 Acq On : 12 Aug 2015 11:13 am
 Operator : TOANP
 Sample : mb
 Misc : MS89345,V4B2296,w,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 12 17:19:20 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	159324	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	450495	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	509837	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	441074	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	215904	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	159782	49.05	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	98.10%
50) 1,2-dichloroethane-d4 (s)	9.02	65	165032	46.10	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	92.20%
80) toluene-d8 (s)	11.02	98	589342	49.10	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	98.20%
105) 4-bromofluorobenzene (s)	13.86	95	197125	52.30	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.60%

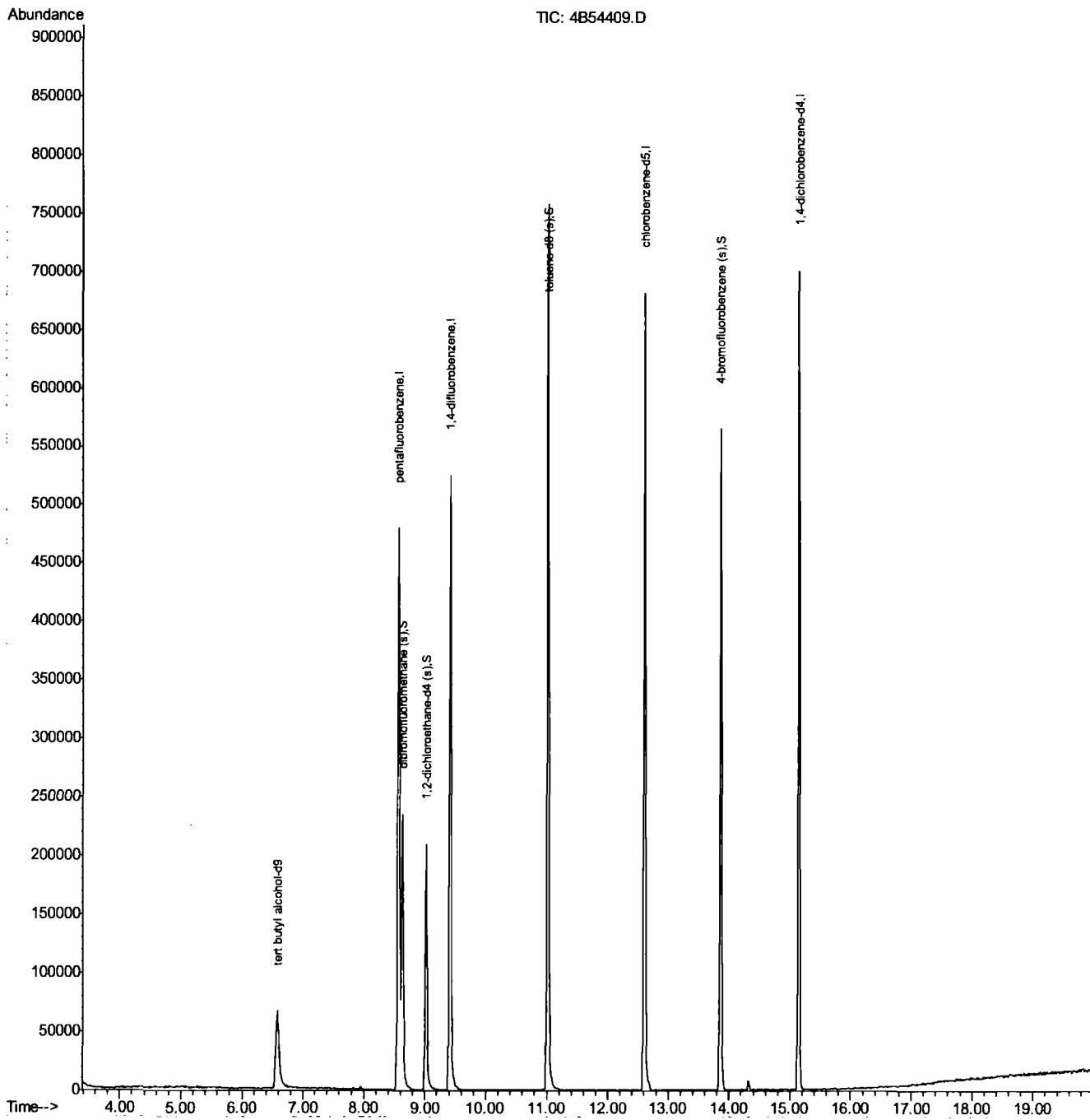
Target Compounds	Qvalue
(#)	= qualifier out of range (m) = manual integration (+) = signals summed

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54409.D
 Acq On : 12 Aug 2015 11:13 am
 Operator : TOANP
 Sample : mb
 Misc : MS89345,V4B2296,w,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 12 17:19:20 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54433.D
 Acq On : 12 Aug 2015 10:30 pm
 Operator : TOANP
 Sample : mb
 Misc : MS89344, V4B2297, w, , , 1
 ALS Vial : 29 Sample Multiplier: 1

Quant Time: Aug 13 12:18:03 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	149010	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	437268	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	487916	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	429546	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	212618	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	155085	49.05	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	98.10%
50) 1,2-dichloroethane-d4 (s)	9.02	65	160972	46.32	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	92.64%
80) toluene-d8 (s)	11.02	98	562290	48.95	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	97.90%
105) 4-bromofluorobenzene (s)	13.86	95	192907	51.97	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	103.94%

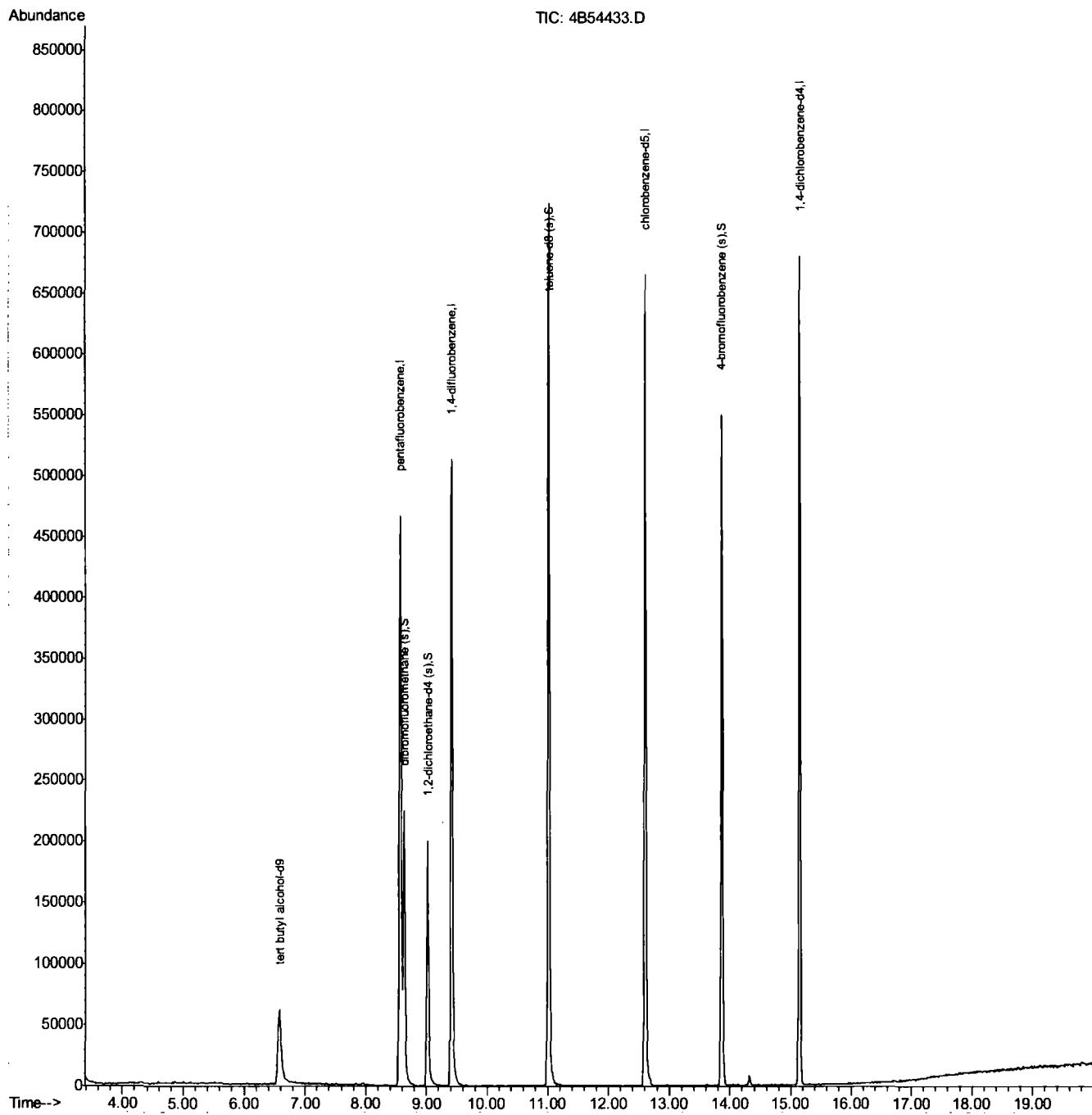
Target Compounds	Qvalue
(#)	= qualifier out of range (m) = manual integration (+) = signals summed

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54433.D
 Acq On : 12 Aug 2015 10:30 pm
 Operator : TOANP
 Sample : mb
 Misc : MS89344,V4B2297,w,,,1
 ALS Vial : 29 Sample Multiplier: 1

Quant Time: Aug 13 12:18:03 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54457.D
 Acq On : 13 Aug 2015 10:43 am
 Operator : TOANP
 Sample : mb2
 Misc : MS89470,V4B2297,w,,,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Aug 13 16:13:50 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	129837	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	435467	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	485700	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	421308	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	200871	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	151604	48.15	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 96.30%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	159509	46.09	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 92.18%	
80) toluene-d8 (s)	11.01	98	564925	49.41	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 98.82%	
105) 4-bromofluorobenzene (s)	13.86	95	186401	53.15	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 106.30%	

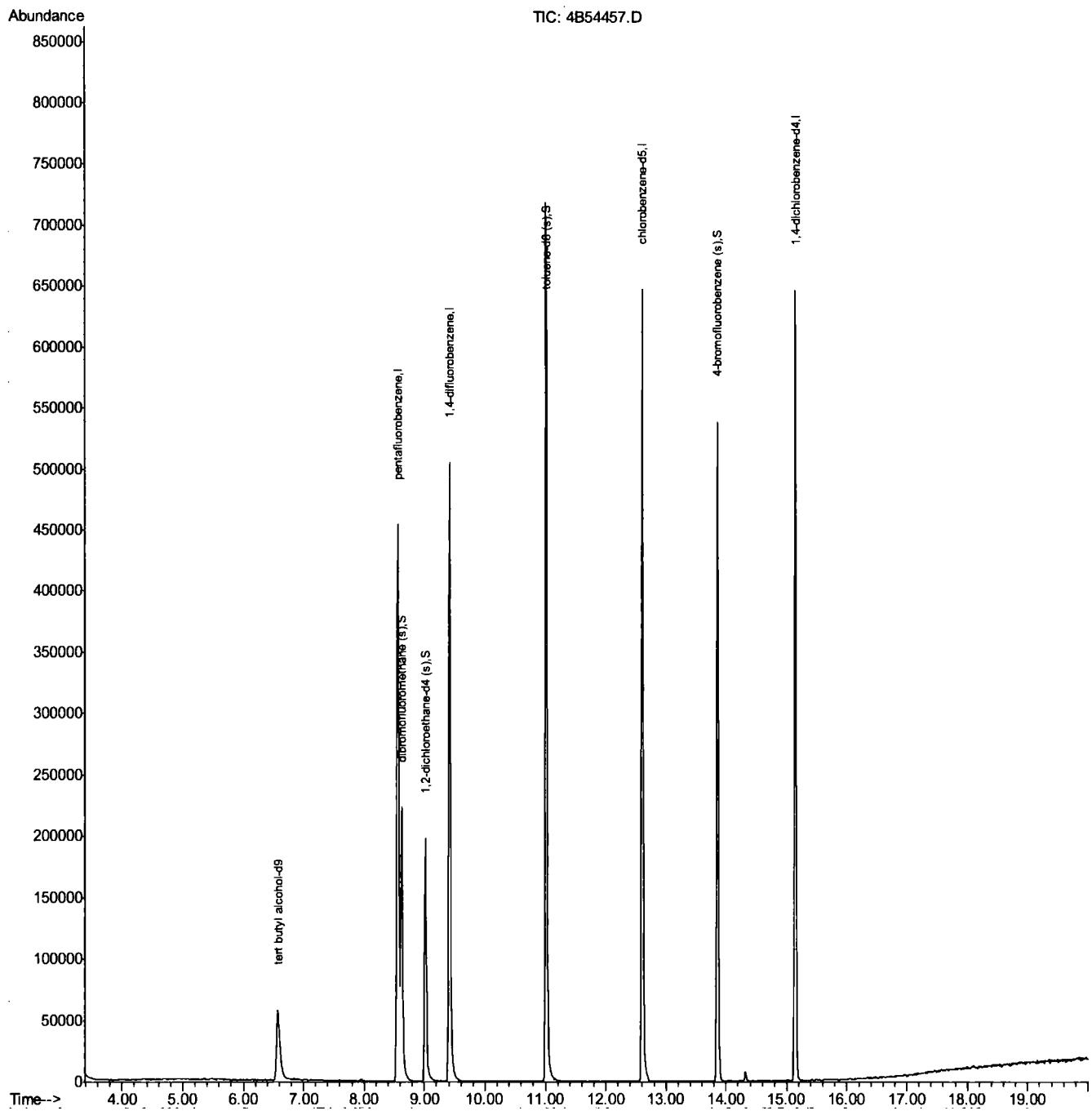
Target Compounds	Qvalue
(#)	= qualifier out of range (m) = manual integration (+) = signals summed

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54457.D
 Acq On : 13 Aug 2015 10:43 am
 Operator : TOANP
 Sample : mb2
 Misc : MS89470,V4B2297,w,,,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Aug 13 16:13:50 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54410.D
 Acq On : 12 Aug 2015 11:47 am
 Operator : TOANP
 Sample : bs
 Misc : MS89342,V4B2296,w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 12 17:19:40 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	125516	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	422985	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	486014	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	432148	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	252161	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	152320	49.80	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 99.60%	
50) 1,2-dichloroethane-d4 (s)	9.01	65	157774	46.94	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 93.88%	
80) toluene-d8 (s)	11.02	98	569670	49.79	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 99.58%	
105) 4-bromofluorobenzene (s)	13.86	95	210611	47.84	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 95.68%	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) tertiary butyl alcohol	6.67	59	80738	265.92	ug/L	92
3) 1,4-dioxane	10.11	88	40983	1460.06	ug/L	96
7) chlorodifluoromethane	3.63	51	208473	48.74	ug/L	97
8) dichlorodifluoromethane	3.60	85	214457	52.35	ug/L	99
10) chloromethane	3.97	52	70440	44.96	ug/L	96
11) vinyl chloride	4.17	62	197387	43.84	ug/L	99
12) bromomethane	4.76	94	139164	53.90	ug/L	97
13) chloroethane	4.91	64	105680	59.70	ug/L	93
15) trichlorofluoromethane	5.28	101	230436	54.19	ug/L	97
19) ethyl ether	5.63	74	75972	47.13	ug/L	93
20) 2-chloropropane	5.81	39	19076	37.19	ug/L	89
21) acrolein	5.86	56	241166	446.86	ug/L	100
22) 1,1-dichloroethene	6.03	96	170653	51.10	ug/L	94
23) acetone	6.05	58	11987	42.66	ug/L	# 68
24) allyl chloride	6.46	76	133255	54.50	ug/L	99
25) acetonitrile	6.44	40	121339	451.09	ug/L	# 65
26) iodomethane	6.28	142	308239	48.87	ug/L	99
27) carbon disulfide	6.38	76	503357	51.36	ug/L	99
28) methylene chloride	6.65	84	175943	49.73	ug/L	95
29) methyl acetate	6.44	74	22821	47.93	ug/L	# 76
30) 1-chloropropane	6.66	42	264519	43.81	ug/L	98
31) methyl tert butyl ether	6.89	73	434034	48.70	ug/L	99
32) trans-1,2-dichloroethene	6.95	96	149297	45.45	ug/L	95
33) di-isopropyl ether	7.41	45	518284	44.35	ug/L	79
34) 2-butanone	8.08	72	17316	46.25	ug/L	# 91
35) 1,1-dichloroethane	7.45	63	288399	46.72	ug/L	100
36) chloroprene	7.54	53	216192	47.29	ug/L	97
37) acrylonitrile	6.92	53	280875	223.42	ug/L	99
38) vinyl acetate	7.43	86	23520	49.32	ug/L	56
39) ethyl tert-butyl ether	7.83	59	503173	48.46	ug/L	99
40) ethyl acetate	8.08	45	21140	41.60	ug/L	92
41) 2,2-dichloropropane	8.11	77	114135	43.73	ug/L	94
42) cis-1,2-dichloroethene	8.11	96	170950	47.91	ug/L	91

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54410.D
 Acq On : 12 Aug 2015 11:47 am
 Operator : TOANP
 Sample : bs
 Misc : MS89342,V4B2296,w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 12 17:19:40 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
43) methylacrylate	8.16	85	23445	48.38	ug/L	# 85
44) propionitrile	8.18	54	212359	418.80	ug/L	93
45) bromochloromethane	8.40	128	92879	50.71	ug/L	92
46) tetrahydrofuran	8.43	42	49194	40.37	ug/L	97
47) chloroform	8.44	85	184819	50.04	ug/L	96
48) T-BUTYL FORMATE	8.46	59	119362	50.56	ug/L	97
51) freon 113	5.97	151	113766	56.11	ug/L	94
52) methacrylonitrile	8.34	41	94903	42.38	ug/L	94
53) 1,1,1-trichloroethane	8.67	97	193415	49.53	ug/L	98
54) cyclohexane	8.74	84	193475	45.21	ug/L	# 80
55) iso-butyl alcohol	8.82	43	62167	446.09	ug/L	96
57) epichlorohydrin	10.63	57	76150	228.23	ug/L	98
58) n-butyl alcohol	9.53	56	227462	2404.11	ug/L	95
59) carbon tetrachloride	8.86	117	192962	49.62	ug/L	98
60) 1,1-dichloropropene	8.83	75	216624	50.13	ug/L	98
61) hexane	7.20	57	163564	49.84	ug/L	96
63) benzene	9.08	78	611879	48.11	ug/L	99
64) iso-octane	9.06	57	469419	45.95	ug/L	98
65) tert-amyl methyl ether	9.09	87	101319	50.11	ug/L	96
66) heptane	9.21	57	140100	62.00	ug/L	95
67) isopropyl acetate	8.99	61	62825	41.32	ug/L	# 89
68) 1,2-dichloroethane	9.10	62	208216	49.68	ug/L	96
69) trichloroethene	9.74	95	174303	50.92	ug/L	99
72) 2-nitropropane	10.50	41	62908	45.85	ug/L	# 79
73) 2-chloroethyl vinyl ether	10.50	63	446068	220.00	ug/L	97
74) methyl methacrylate	9.99	100	43054	52.78	ug/L	97
75) 1,2-dichloropropane	10.00	63	174236	47.69	ug/L	98
76) dibromomethane	10.16	93	107104	50.34	ug/L	96
77) methylcyclohexane	9.94	83	239493	54.41	ug/L	95
78) bromodichloromethane	10.27	83	233335	51.00	ug/L	98
79) cis-1,3-dichloropropene	10.72	75	283004	47.74	ug/L	97
81) 4-methyl-2-pentanone	10.82	58	66400	47.28	ug/L	98
82) toluene	11.09	92	393840	48.58	ug/L	98
83) 3-methyl-1-butanol	10.83	55	140042	893.14	ug/L	94
84) trans-1,3-dichloropropene	11.29	75	244719	47.14	ug/L	98
85) ethyl methacrylate	11.27	69	208579	48.56	ug/L	95
86) 1,1,2-trichloroethane	11.52	83	135123	47.88	ug/L	97
87) 2-hexanone	11.69	58	58096	45.30	ug/L	89
89) tetrachloroethene	11.69	164	190106	50.64	ug/L	100
90) 1,3-dichloropropane	11.71	76	243801	47.93	ug/L	95
91) butyl acetate	11.76	56	102078	44.00	ug/L	97
92) 3,3-DIMETHYL-1-BUTANOL	11.86	57	157177	461.68	ug/L	97
93) dibromochloromethane	11.99	129	200308	50.36	ug/L	98
94) 1,2-dibromoethane	12.15	107	169586	50.90	ug/L	98
96) chlorobenzene	12.63	112	462805	50.64	ug/L	97
97) 1,1,1,2-tetrachloroethane	12.69	131	171668	49.15	ug/L	98
98) ethylbenzene	12.68	91	759642	49.90	ug/L	98
99) m,p-xylene	12.80	106	588889	101.52	ug/L	98
100) o-xylene	13.25	106	309374	52.90	ug/L	99
101) styrene	13.27	104	504694	50.54	ug/L	96

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54410.D
 Acq On : 12 Aug 2015 11:47 am
 Operator : TOANP
 Sample : bs
 Misc : MS89342,V4B2296,w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 12 17:19:40 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

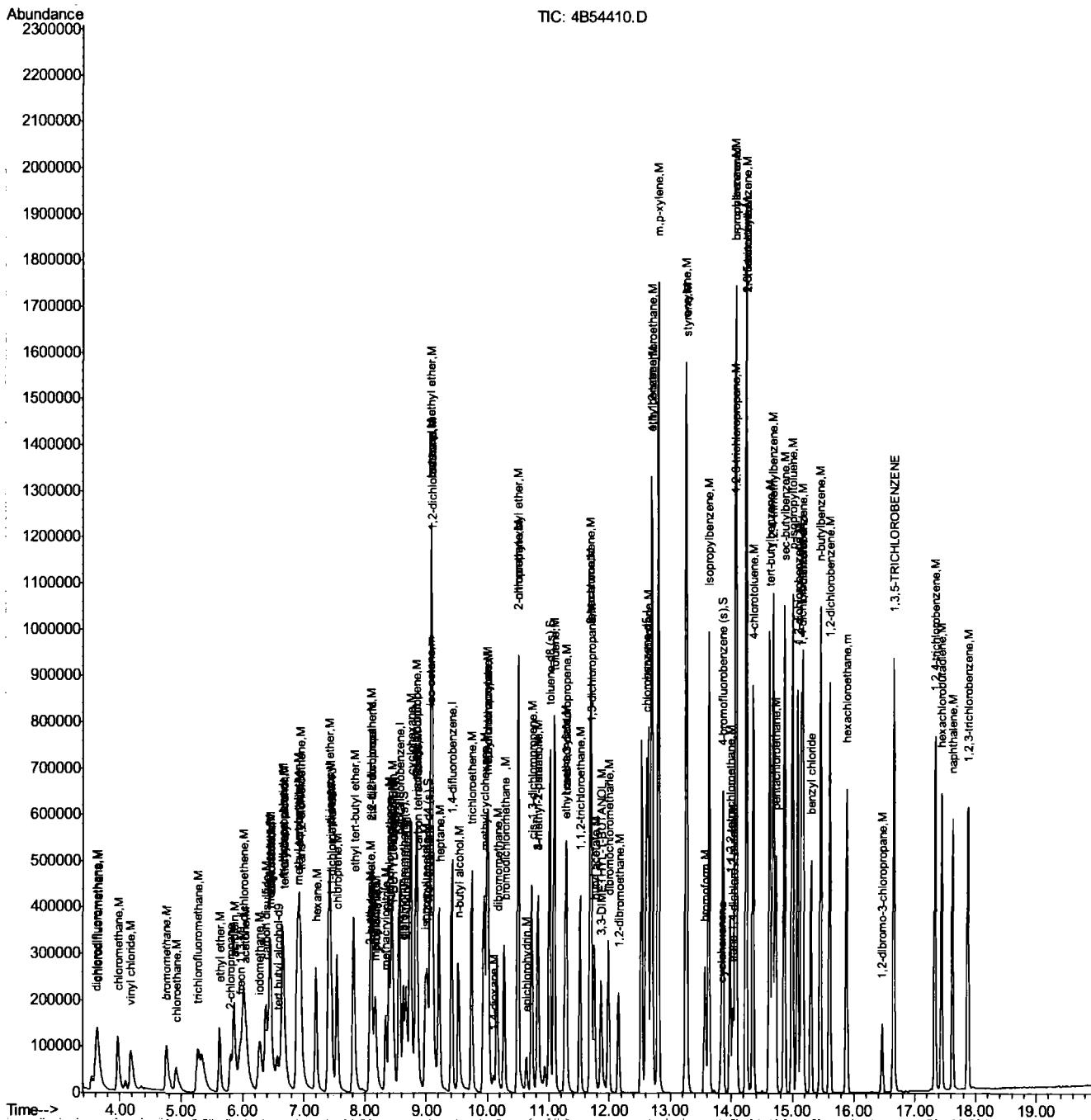
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
102) bromoform	13.57	173	154764	50.07	ug/L	97
104) isopropylbenzene	13.63	105	773988	48.35	ug/L	99
106) cyclohexanone	13.82	55	58265	127.57	ug/L	98
107) bromobenzene	14.07	156	234879	50.02	ug/L	94
108) 1,1,2,2-tetrachloroethane	13.98	83	220431	47.02	ug/L	99
109) trans-1,4-dichloro-2-buten	14.02	53	40192	33.35	ug/L	98
110) 1,2,3-trichloropropane	14.05	110	54350	48.93	ug/L	97
111) n-propylbenzene	14.08	91	913646	50.49	ug/L	98
113) 2-chlorotoluene	14.24	126	196888	48.01	ug/L	95
114) 4-chlorotoluene	14.35	91	559301	47.81	ug/L	100
115) 1,3,5-trimethylbenzene	14.24	105	651161	48.32	ug/L	99
116) tert-butylbenzene	14.62	119	593009	50.73	ug/L	99
117) pentachloroethane	14.72	167	133059	55.00	ug/L	97
118) 1,2,4-trimethylbenzene	14.68	105	670628	51.67	ug/L	99
119) sec-butylbenzene	14.87	105	865965	48.76	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	430996	50.43	ug/L	99
121) p-isopropyltoluene	15.00	119	747967	50.37	ug/L	98
122) 1,4-dichlorobenzene	15.17	146	429954	50.29	ug/L	99
123) benzyl chloride	15.30	91	406259	47.55	ug/L	99
124) 1,2-dichlorobenzene	15.61	146	432194	51.78	ug/L	98
126) n-butylbenzene	15.46	92	369492	50.17	ug/L	98
128) 1,2-dibromo-3-chloropropan	16.46	75	37125	49.34	ug/L	98
129) 1,3,5-TRICHLOROBENZENE	16.65	180	359822	54.06	ug/L	99
130) 1,2,4-trichlorobenzene	17.33	180	277779	52.10	ug/L	99
131) hexachlorobutadiene	17.44	225	172131	53.05	ug/L	98
132) naphthalene	17.62	128	486651	51.92	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	239182	53.19	ug/L	99
134) hexachloroethane	15.89	201	157654	51.21	ug/L	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54410.D
Acq On : 12 Aug 2015 11:47 am
Operator : TOANP
Sample : bs
Misc : MS89342,V4B2296,w,,,1
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 12 17:19:40 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54434.D
 Acq On : 12 Aug 2015 10:58 pm
 Operator : TOANP
 Sample : bs
 Misc : MS89499,V4B2297,w,,,1
 ALS Vial : 30 Sample Multiplier: 1

Quant Time: Aug 13 12:18:14 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.57	65	119545	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	416714	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	479398	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	432369	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	251910	50.00	ug/L	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
49) dibromofluoromethane (s)	8.63	113	150569	49.97	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery = 99.94%			
50) 1,2-dichloroethane-d4 (s)	9.02	65	156539	47.27	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery = 94.54%			
80) toluene-d8 (s)	11.02	98	562727	49.86	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery = 99.72%			
105) 4-bromofluorobenzene (s)	13.86	95	211529	48.10	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 96.20%			

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) tertiary butyl alcohol	6.67	59	77819	269.11	ug/L	81
3) 1,4-dioxane	10.12	88	39076	1461.65	ug/L	90
7) chlorodifluoromethane	3.64	51	194732	46.21	ug/L	96
8) dichlorodifluoromethane	3.61	85	209439	51.90	ug/L	97
10) chloromethane	3.96	52	63419	41.09	ug/L	95
11) vinyl chloride	4.18	62	190998	43.06	ug/L	98
12) bromomethane	4.76	94	125552	49.36	ug/L	99
13) chloroethane	4.91	64	97508	55.91	ug/L	95
15) trichlorofluoromethane	5.28	101	218388	52.13	ug/L	96
19) ethyl ether	5.63	74	74125	46.68	ug/L	93
20) 2-chloropropane	5.81	39	27738	54.89	ug/L	72
21) acrolein	5.86	56	230474	433.47	ug/L	98
22) 1,1-dichloroethene	6.02	96	158620	48.21	ug/L	95
23) acetone	6.05	58	12338	44.57	ug/L	100
24) allyl chloride	6.46	76	92133	38.25	ug/L	# 69
25) acetonitrile	6.44	40	111822	421.97	ug/L	# 61
26) iodomethane	6.28	142	318367	51.24	ug/L	98
27) carbon disulfide	6.38	76	496936	51.46	ug/L	100
28) methylene chloride	6.64	84	167070	47.93	ug/L	89
29) methyl acetate	6.44	74	22655	48.29	ug/L	92
30) 1-chloropropane	6.66	42	252662	42.48	ug/L	98
31) methyl tert butyl ether	6.89	73	420931	47.95	ug/L	97
32) trans-1,2-dichloroethene	6.95	96	148779	45.97	ug/L	94
33) di-isopropyl ether	7.41	45	503737	43.75	ug/L	83
34) 2-butanone	8.08	72	17274	46.84	ug/L	# 71
35) 1,1-dichloroethane	7.45	63	277889	45.69	ug/L	99
36) chloroprene	7.54	53	215684	47.89	ug/L	99
37) acrylonitrile	6.92	53	275960	222.81	ug/L	99
38) vinyl acetate	7.43	86	22768	48.46	ug/L	94
39) ethyl tert-butyl ether	7.83	59	484932	47.41	ug/L	99
40) ethyl acetate	8.08	45	21146	42.24	ug/L	# 86
41) 2,2-dichloropropane	8.11	77	105539	41.05	ug/L	96
42) cis-1,2-dichloroethene	8.11	96	166008	47.22	ug/L	92

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54434.D
 Acq On : 12 Aug 2015 10:58 pm
 Operator : TOANP
 Sample : bs
 Misc : MS89499,V4B2297,w,,,1
 ALS Vial : 30 Sample Multiplier: 1

Quant Time: Aug 13 12:18:14 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.16	85	23370	48.95	ug/L	# 83
44) propionitrile	8.18	54	205034	410.44	ug/L	99
45) bromochloromethane	8.40	128	90308	50.05	ug/L	88
46) tetrahydrofuran	8.43	42	48682	40.55	ug/L	94
47) chloroform	8.44	85	176689	48.56	ug/L	99
48) T-BUTYL FORMATE	8.46	59	117095	50.34	ug/L	98
51) freon 113	5.97	151	109688	54.91	ug/L	93
52) methacrylonitrile	8.34	41	92804	42.06	ug/L	94
53) 1,1,1-trichloroethane	8.67	97	200468	52.10	ug/L	99
54) cyclohexane	8.74	84	195425	46.35	ug/L	98
55) iso-butyl alcohol	8.82	43	52898	385.29	ug/L	96
57) epichlorohydrin	10.63	57	70926	215.51	ug/L	97
58) n-butyl alcohol	9.53	56	220654	2364.34	ug/L	95
59) carbon tetrachloride	8.86	117	203113	52.95	ug/L	96
60) 1,1-dichloropropene	8.83	75	207673	48.73	ug/L	98
61) hexane	7.20	57	147470	45.55	ug/L	96
63) benzene	9.08	78	599187	47.76	ug/L	99
64) iso-octane	9.07	57	424347	42.11	ug/L	99
65) tert-amyl methyl ether	9.09	87	100064	50.17	ug/L	91
66) heptane	9.21	57	122580	55.00	ug/L	95
67) isopropyl acetate	8.99	61	62596	41.74	ug/L	96
68) 1,2-dichloroethane	9.10	62	206975	50.07	ug/L	97
69) trichloroethene	9.74	95	170107	50.38	ug/L	98
72) 2-nitropropane	10.50	41	63610	47.00	ug/L	# 75
73) 2-chloroethyl vinyl ether	10.50	63	451101	225.55	ug/L	97
74) methyl methacrylate	9.99	100	43032	53.48	ug/L	94
75) 1,2-dichloropropane	10.00	63	169332	46.98	ug/L	100
76) dibromomethane	10.16	93	105542	50.29	ug/L	96
77) methylcyclohexane	9.94	83	228391	52.60	ug/L	96
78) bromodichloromethane	10.28	83	230393	51.05	ug/L	96
79) cis-1,3-dichloropropene	10.72	75	269693	46.12	ug/L	99
81) 4-methyl-2-pentanone	10.82	58	65511	47.29	ug/L	99
82) toluene	11.09	92	389391	48.70	ug/L	100
83) 3-methyl-1-butanol	10.83	55	135351	875.14	ug/L	94
84) trans-1,3-dichloropropene	11.29	75	238087	46.49	ug/L	97
85) ethyl methacrylate	11.27	69	208989	49.33	ug/L	94
86) 1,1,2-trichloroethane	11.52	83	133047	47.80	ug/L	100
87) 2-hexanone	11.69	58	57589	45.53	ug/L	97
89) tetrachloroethene	11.69	164	190649	50.76	ug/L	99
90) 1,3-dichloropropane	11.71	76	240641	47.29	ug/L	98
91) butyl acetate	11.76	56	100090	43.12	ug/L	97
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	148539	436.09	ug/L	98
93) dibromochloromethane	11.99	129	198751	49.95	ug/L	98
94) 1,2-dibromoethane	12.15	107	170201	51.06	ug/L	99
96) chlorobenzene	12.63	112	461919	50.51	ug/L	97
97) 1,1,1,2-tetrachloroethane	12.70	131	171673	49.12	ug/L	97
98) ethylbenzene	12.68	91	747421	49.07	ug/L	98
99) m,p-xylene	12.80	106	577161	99.44	ug/L	97
100) o-xylene	13.25	106	301613	51.55	ug/L	98
101) styrene	13.27	104	499698	50.01	ug/L	97

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54434.D
 Acq On : 12 Aug 2015 10:58 pm
 Operator : TOANP
 Sample : bs
 Misc : MS89499,V4B2297,w,,,1
 ALS Vial : 30 Sample Multiplier: 1

Quant Time: Aug 13 12:18:14 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

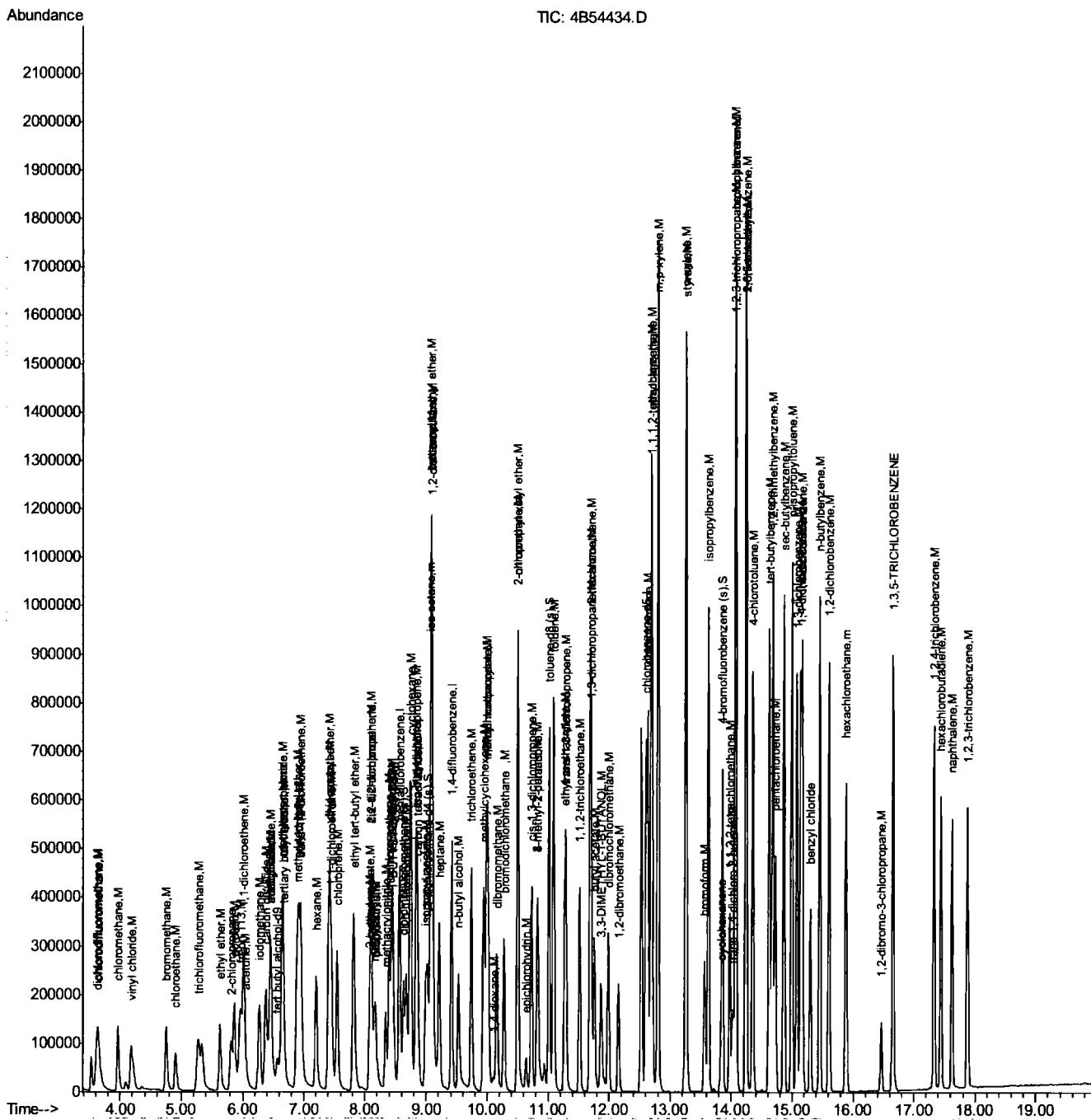
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
102) bromoform	13.57	173	154475	49.95	ug/L	97
104) isopropylbenzene	13.63	105	757885	47.39	ug/L	99
106) cyclohexanone	13.83	55	51567	113.02	ug/L	97
107) bromobenzene	14.07	156	234384	49.97	ug/L	94
108) 1,1,2,2-tetrachloroethane	13.98	83	216775	46.29	ug/L	100
109) trans-1,4-dichloro-2-buten	14.02	53	36579	30.38	ug/L	96
110) 1,2,3-trichloropropane	14.06	110	54601	49.21	ug/L	94
111) n-propylbenzene	14.08	91	900669	49.83	ug/L	98
113) 2-chlorotoluene	14.24	126	194849	47.56	ug/L	94
114) 4-chlorotoluene	14.35	91	549230	46.99	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	638965	47.47	ug/L	98
116) tert-butylbenzene	14.62	119	579694	49.64	ug/L	98
117) pentachloroethane	14.72	167	124082	51.34	ug/L	96
118) 1,2,4-trimethylbenzene	14.68	105	663282	51.16	ug/L	99
119) sec-butylbenzene	14.87	105	842420	47.48	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	426355	49.93	ug/L	99
121) p-isopropyltoluene	15.00	119	733198	49.43	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	421366	49.34	ug/L	99
123) benzyl chloride	15.30	91	303218	35.52	ug/L	98
124) 1,2-dichlorobenzene	15.61	146	425305	51.01	ug/L	98
126) n-butylbenzene	15.46	92	353705	48.08	ug/L	99
128) 1,2-dibromo-3-chloropropan	16.46	75	35318	46.99	ug/L	94
129) 1,3,5-TRICHLOROBENZENE	16.65	180	349714	52.59	ug/L	99
130) 1,2,4-trichlorobenzene	17.33	180	268201	50.35	ug/L	99
131) hexachlorobutadiene	17.44	225	165651	51.10	ug/L	98
132) naphthalene	17.62	128	469696	50.16	ug/L	100
133) 1,2,3-trichlorobenzene	17.88	180	227814	50.71	ug/L	99
134) hexachloroethane	15.89	201	154070	50.09	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (OT Reviewed)

Data Path : C:\MSDChem\1\DATA\
Data File : 4B54434.D
Acq On : 12 Aug 2015 10:58 pm
Operator : TOANP
Sample : bs
Misc : MS89499,V4B2297,w,,,1
ALS Vial : 30 Sample Multiplier: 1

Quant Time: Aug 13 12:18:14 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54420.D
 Acq On : 12 Aug 2015 4:27 pm
 Operator : TOANP
 Sample : jc1107-3ms
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Aug 12 17:34:23 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	145924	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	413150	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	475209	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	424091	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	251934	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	146219	48.94	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 97.88%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	154947	47.19	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 94.38%	
80) toluene-d8 (s)	11.02	98	548875	49.06	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 98.12%	
105) 4-bromofluorobenzene (s)	13.86	95	209894	47.72	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 95.44%	

Target Compounds

				QValue	
2) tertiary butyl alcohol	6.68	59	92446	261.90	ug/L # 66
3) 1,4-dioxane	10.11	88	46419	1422.44	ug/L 91
7) chlorodifluoromethane	3.64	51	202387	48.44	ug/L 97
8) dichlorodifluoromethane	3.62	85	182817	45.69	ug/L 99
10) chloromethane	3.96	52	59267	38.73	ug/L 95
11) vinyl chloride	4.17	62	185509	42.19	ug/L 96
12) bromomethane	4.75	94	116429	46.17	ug/L 96
13) chloroethane	4.91	64	80362	46.48	ug/L 96
15) trichlorofluoromethane	5.28	101	229438	55.24	ug/L 93
19) ethyl ether	5.63	74	77152	49.00	ug/L 95
20) 2-chloropropane	5.81	39	28732	57.35	ug/L 83
21) acrolein	5.86	56	169159	320.90	ug/L 97
22) 1,1-dichloroethene	6.02	96	170664	52.32	ug/L 93
23) acetone	6.06	58	11899	43.36	ug/L 87
24) allyl chloride	6.46	76	104754	43.86	ug/L # 73
25) acetonitrile	6.45	40	121562	462.68	ug/L # 77
26) iodomethane	6.27	142	333529	54.14	ug/L 100
27) carbon disulfide	6.38	76	515813	53.88	ug/L 99
28) methylene chloride	6.64	84	173444	50.19	ug/L 88
29) methyl acetate	6.44	74	24287	52.22	ug/L # 82
30) 1-chloropropane	6.66	42	273886	46.44	ug/L 98
31) methyl tert butyl ether	6.89	73	444395	51.05	ug/L 97
32) trans-1,2-dichloroethene	6.95	96	159396	49.68	ug/L 94
33) di-isopropyl ether	7.41	45	509304	44.62	ug/L 85
34) 2-butanone	8.08	72	18189	49.74	ug/L # 92
35) 1,1-dichloroethane	7.46	63	304212	50.45	ug/L 98
36) chloroprene	7.55	53	223363	50.02	ug/L 96
37) acrylonitrile	6.92	53	284746	231.89	ug/L 99
38) vinyl acetate	7.43	86	25221	54.14	ug/L 91
39) ethyl tert-butyl ether	7.83	59	492874	48.60	ug/L 98
40) ethyl acetate	8.08	45	22167	44.66	ug/L # 82
41) 2,2-dichloropropane	8.11	77	132587	52.01	ug/L 98
42) cis-1,2-dichloroethene	8.11	96	175506	50.36	ug/L 92

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54420.D
 Acq On : 12 Aug 2015 4:27 pm
 Operator : TOANP
 Sample : jc1107-3ms
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Aug 12 17:34:23 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
43) methylacrylate	8.16	85	24522	51.81	ug/L	# 90
44) propionitrile	8.18	54	235959	476.42	ug/L	98
45) bromochloromethane	8.40	128	96004	53.67	ug/L	88
46) tetrahydrofuran	8.43	42	49313	41.43	ug/L	95
47) chloroform	8.44	85	187825	52.06	ug/L	96
48) T-BUTYL FORMATE	8.47	59	69522	30.15	ug/L	92
51) freon 113	5.97	151	118220	59.69	ug/L	89
52) methacrylonitrile	8.34	41	97714	44.67	ug/L	93
53) 1,1,1-trichloroethane	8.67	97	223758	58.66	ug/L	98
54) cyclohexane	8.74	84	226065	54.08	ug/L	86
55) iso-butyl alcohol	8.82	43	64461	473.57	ug/L	94
57) epichlorohydrin	10.63	57	74823	229.35	ug/L	96
58) n-butyl alcohol	9.53	56	264824	2862.64	ug/L	96
59) carbon tetrachloride	8.86	117	215871	56.77	ug/L	99
60) 1,1-dichloropropene	8.84	75	219639	51.99	ug/L	97
61) hexane	7.20	57	184761	57.58	ug/L	95
63) benzene	9.08	78	620517	49.90	ug/L	100
64) iso-octane	9.07	57	522751	52.33	ug/L	92
65) tert-amyl methyl ether	9.09	87	99789	50.47	ug/L	94
66) heptane	9.21	57	146747	66.42	ug/L	97
67) isopropyl acetate	8.99	61	64892	43.65	ug/L	93
68) 1,2-dichloroethane	9.10	62	214202	52.27	ug/L	96
69) trichloroethene	9.74	95	179466	53.62	ug/L	98
72) 2-nitropropane	10.50	41	48292	35.99	ug/L	# 1
74) methyl methacrylate	9.99	100	44117	55.31	ug/L	95
75) 1,2-dichloropropane	10.00	63	170253	47.66	ug/L	99
76) dibromomethane	10.16	93	111304	53.50	ug/L	96
77) methylcyclohexane	9.94	83	252717	58.72	ug/L	97
78) bromodichloromethane	10.28	83	239082	53.44	ug/L	97
79) cis-1,3-dichloropropene	10.72	75	290900	50.19	ug/L	99
81) 4-methyl-2-pentanone	10.82	58	70838	51.59	ug/L	100
82) toluene	11.09	92	407614	51.42	ug/L	100
83) 3-methyl-1-butanol	10.83	55	170672	1113.24	ug/L	93
84) trans-1,3-dichloropropene	11.29	75	263522	51.91	ug/L	98
85) ethyl methacrylate	11.27	69	224209	53.39	ug/L	95
86) 1,1,2-trichloroethane	11.52	83	139070	50.40	ug/L	99
87) 2-hexanone	11.69	58	63467	50.62	ug/L	91
89) tetrachloroethene	11.69	164	199308	54.10	ug/L	99
90) 1,3-dichloropropane	11.71	76	256067	51.30	ug/L	96
91) butyl acetate	11.76	56	110480	48.52	ug/L	90
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	205818	616.04	ug/L	98
93) dibromochloromethane	11.99	129	208162	53.33	ug/L	100
94) 1,2-dibromoethane	12.15	107	178685	54.65	ug/L	100
96) chlorobenzene	12.63	112	480473	53.57	ug/L	99
97) 1,1,1,2-tetrachloroethane	12.69	131	184198	53.73	ug/L	99
98) ethylbenzene	12.68	91	794714	53.19	ug/L	98
99) m,p-xylene	12.80	106	615541	108.13	ug/L	98
100) o-xylene	13.25	106	315041	54.89	ug/L	98
101) styrene	13.27	104	529958	54.07	ug/L	96
102) bromoform	13.57	173	162017	53.41	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54420.D
 Acq On : 12 Aug 2015 4:27 pm
 Operator : TOANP
 Sample : jc1107-3ms
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 16 Sample Multiplier: 1

Quant Time: Aug 12 17:34:23 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

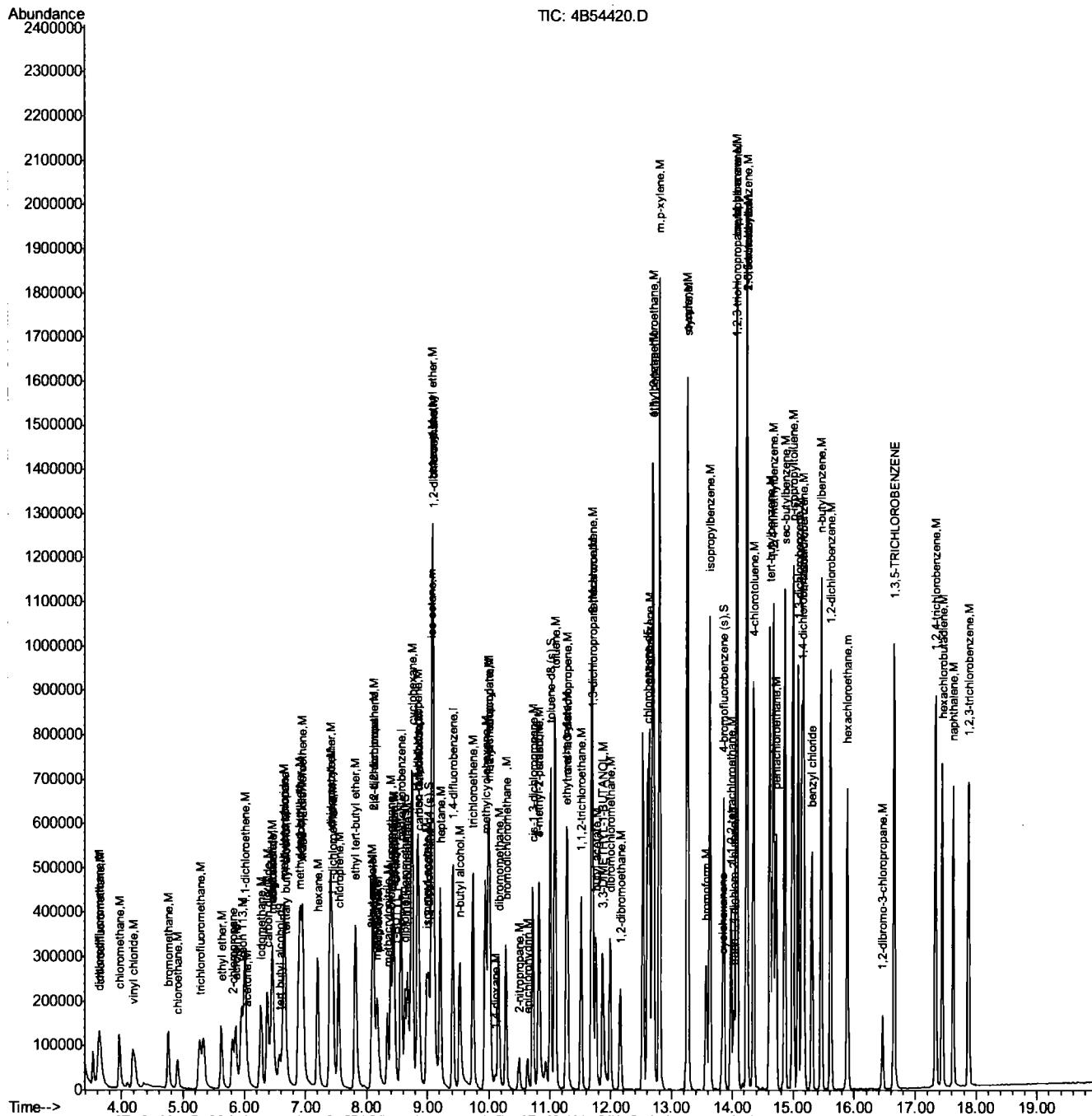
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
104) isopropylbenzene	13.63	105	815736	51.01	ug/L	99
106) cyclohexanone	13.83	55	66806	146.40	ug/L	96
107) bromobenzene	14.07	156	248024	52.87	ug/L	94
108) 1,1,2,2-tetrachloroethane	13.98	83	246827	52.70	ug/L	97
109) trans-1,4-dichloro-2-buten	14.02	53	36786	30.55	ug/L	98
110) 1,2,3-trichloropropane	14.06	110	59661	53.76	ug/L	97
111) n-propylbenzene	14.08	91	923668	51.09	ug/L	98
113) 2-chlorotoluene	14.24	126	209162	51.05	ug/L	93
114) 4-chlorotoluene	14.35	91	591958	50.64	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	694141	51.56	ug/L	99
116) tert-butylbenzene	14.62	119	623349	53.38	ug/L	100
117) pentachloroethane	14.72	167	145768	60.31	ug/L	97
118) 1,2,4-trimethylbenzene	14.68	105	688126	53.07	ug/L	99
119) sec-butylbenzene	14.87	105	936467	52.77	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	467336	54.73	ug/L	99
121) p-isopropyltoluene	15.00	119	797547	53.76	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	458472	53.68	ug/L	99
123) benzyl chloride	15.30	91	442896	51.88	ug/L	99
124) 1,2-dichlorobenzene	15.61	146	460120	55.18	ug/L	99
126) n-butylbenzene	15.46	92	398677	54.19	ug/L	99
128) 1,2-dibromo-3-chloropropan	16.46	75	41192	54.80	ug/L	98
129) 1,3,5-TRICHLOROBENZENE	16.65	180	394491	59.32	ug/L	100
130) 1,2,4-trichlorobenzene	17.33	180	318498	59.79	ug/L	99
131) hexachlorobutadiene	17.44	225	194742	60.07	ug/L	99
132) naphthalene	17.62	128	564473	60.28	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	268620	59.79	ug/L	99
134) hexachloroethane	15.89	201	163993	53.31	ug/L	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54420.D
Acq On : 12 Aug 2015 4:27 pm
Operator : TOANP
Sample : jc1107-3ms
Misc : MS89470, V4B2296,w,,,1
ALS Vial : 16 Sample Multiplier: 1

Quant Time: Aug 12 17:34:23 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54421.D
 Acq On : 12 Aug 2015 4:55 pm
 Operator : TOANP
 Sample : jc1107-3msd
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Aug 12 17:34:44 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	152325	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	431587	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	496307	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	441468	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	260200	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	154857	49.62	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 99.24%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	161657	47.13	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 94.26%	
80) toluene-d8 (s)	11.02	98	576789	49.37	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 98.74%	
105) 4-bromofluorobenzene (s)	13.86	95	216499	47.66	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 95.32%	

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.67	59	89567	243.08	ug/L
3) 1,4-dioxane	10.11	88	47520	1394.99	ug/L
7) chlorodifluoromethane	3.64	51	210227	48.17	ug/L
8) dichlorodifluoromethane	3.61	85	189508	45.34	ug/L
10) chloromethane	3.96	52	64753	40.51	ug/L
11) vinyl chloride	4.18	62	197226	42.94	ug/L
12) bromomethane	4.75	94	125764	47.74	ug/L
13) chloroethane	4.91	64	87858	48.64	ug/L
15) trichlorofluoromethane	5.27	101	232716	53.63	ug/L
19) ethyl ether	5.63	74	82251	50.01	ug/L
20) 2-chloropropane	5.81	39	29336	56.05	ug/L
21) acrolein	5.87	56	189985	345.01	ug/L
22) 1,1-dichloroethene	6.02	96	177385	52.05	ug/L
23) acetone	6.06	58	13198	46.04	ug/L
24) allyl chloride	6.46	76	108679	43.56	ug/L #
25) acetonitrile	6.46	40	119799	436.49	ug/L
26) iodomethane	6.27	142	347146	53.95	ug/L
27) carbon disulfide	6.38	76	537414	53.74	ug/L
28) methylene chloride	6.64	84	180316	49.95	ug/L
29) methyl acetate	6.44	74	25461	52.41	ug/L #
30) 1-chloropropane	6.66	42	275918	44.79	ug/L
31) methyl tert butyl ether	6.89	73	462206	50.83	ug/L
32) trans-1,2-dichloroethene	6.95	96	165563	49.40	ug/L
33) di-isopropyl ether	7.41	45	524081	43.95	ug/L
34) 2-butanone	8.08	72	19054	49.88	ug/L #
35) 1,1-dichloroethane	7.45	63	314200	49.88	ug/L
36) chloroprene	7.55	53	229938	49.29	ug/L
37) acrylonitrile	6.92	53	302982	236.20	ug/L
38) vinyl acetate	7.43	86	25152	51.69	ug/L
39) ethyl tert-butyl ether	7.83	59	514175	48.54	ug/L
40) ethyl acetate	8.08	45	24201	46.67	ug/L #
41) 2,2-dichloropropane	8.11	77	134937	50.67	ug/L
42) cis-1,2-dichloroethene	8.11	96	182838	50.22	ug/L

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54421.D
 Acq On : 12 Aug 2015 4:55 pm
 Operator : TOANP
 Sample : jc1107-3msd
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Aug 12 17:34:44 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
43) methylacrylate	8.16	85	25257	51.08	ug/L	# 84
44) propionitrile	8.19	54	241220	466.23	ug/L	89
45) bromochloromethane	8.40	128	100267	53.66	ug/L	89
46) tetrahydrofuran	8.43	42	49924	40.16	ug/L	96
47) chloroform	8.44	85	190700	50.60	ug/L	97
48) T-BUTYL FORMATE	8.47	59	58305	24.20	ug/L	87
51) freon 113	5.97	151	123260	59.58	ug/L	93
52) methacrylonitrile	8.34	41	101860	44.58	ug/L	94
53) 1,1,1-trichloroethane	8.67	97	228132	57.25	ug/L	97
54) cyclohexane	8.74	84	232167	53.17	ug/L	87
55) iso-butyl alcohol	8.83	43	63854	449.07	ug/L	97
57) epichlorohydrin	10.64	57	73240	214.96	ug/L	98
58) n-butyl alcohol	9.53	56	275172	2848.05	ug/L	94
59) carbon tetrachloride	8.86	117	218618	55.05	ug/L	99
60) 1,1-dichloropropene	8.83	75	222772	50.49	ug/L	97
61) hexane	7.20	57	190343	56.79	ug/L	98
63) benzene	9.08	78	644086	49.59	ug/L	99
64) iso-octane	9.06	57	525836	50.40	ug/L	91
65) tert-amyl methyl ether	9.09	87	104909	50.81	ug/L	97
66) heptane	9.21	57	152250	65.98	ug/L	98
67) isopropyl acetate	8.99	61	68520	44.13	ug/L	95
68) 1,2-dichloroethane	9.10	62	218971	51.17	ug/L	97
69) trichloroethene	9.74	95	184187	52.69	ug/L	97
72) 2-nitropropane	10.50	41	48732	34.78	ug/L	# 1
74) methyl methacrylate	9.99	100	46827	56.22	ug/L	93
75) 1,2-dichloropropane	10.00	63	180659	48.42	ug/L	98
76) dibromomethane	10.16	93	114696	52.79	ug/L	95
77) methylcyclohexane	9.94	83	257454	57.27	ug/L	97
78) bromodichloromethane	10.28	83	244662	52.36	ug/L	98
79) cis-1,3-dichloropropene	10.72	75	304397	50.29	ug/L	98
81) 4-methyl-2-pentanone	10.82	58	75678	52.77	ug/L	98
82) toluene	11.09	92	421834	50.96	ug/L	100
83) 3-methyl-1-butanol	10.83	55	171711	1072.40	ug/L	93
84) trans-1,3-dichloropropene	11.29	75	271958	51.30	ug/L	99
85) ethyl methacrylate	11.27	69	236210	53.86	ug/L	94
86) 1,1,2-trichloroethane	11.52	83	146215	50.74	ug/L	97
87) 2-hexanone	11.69	58	67440	51.50	ug/L	92
89) tetrachloroethene	11.69	164	203518	53.07	ug/L	99
90) 1,3-dichloropropane	11.71	76	263415	50.70	ug/L	98
91) butyl acetate	11.76	56	113316	47.81	ug/L	91
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	209509	602.41	ug/L	97
93) dibromochloromethane	11.99	129	215971	53.15	ug/L	100
94) 1,2-dibromoethane	12.15	107	185856	54.61	ug/L	100
96) chlorobenzene	12.63	112	497434	53.28	ug/L	98
97) 1,1,1,2-tetrachloroethane	12.70	131	186022	52.13	ug/L	99
98) ethylbenzene	12.69	91	817250	52.55	ug/L	98
99) m,p-xylene	12.80	106	630210	106.35	ug/L	99
100) o-xylene	13.25	106	326053	54.57	ug/L	97
101) styrene	13.27	104	546084	53.53	ug/L	96
102) bromoform	13.57	173	167897	53.17	ug/L	99

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54421.D
 Acq On : 12 Aug 2015 4:55 pm
 Operator : TOANP
 Sample : jc1107-3msd
 Misc : MS89470,V4B2296,w,,,1
 ALS Vial : 17 Sample Multiplier: 1

Quant Time: Aug 12 17:34:44 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

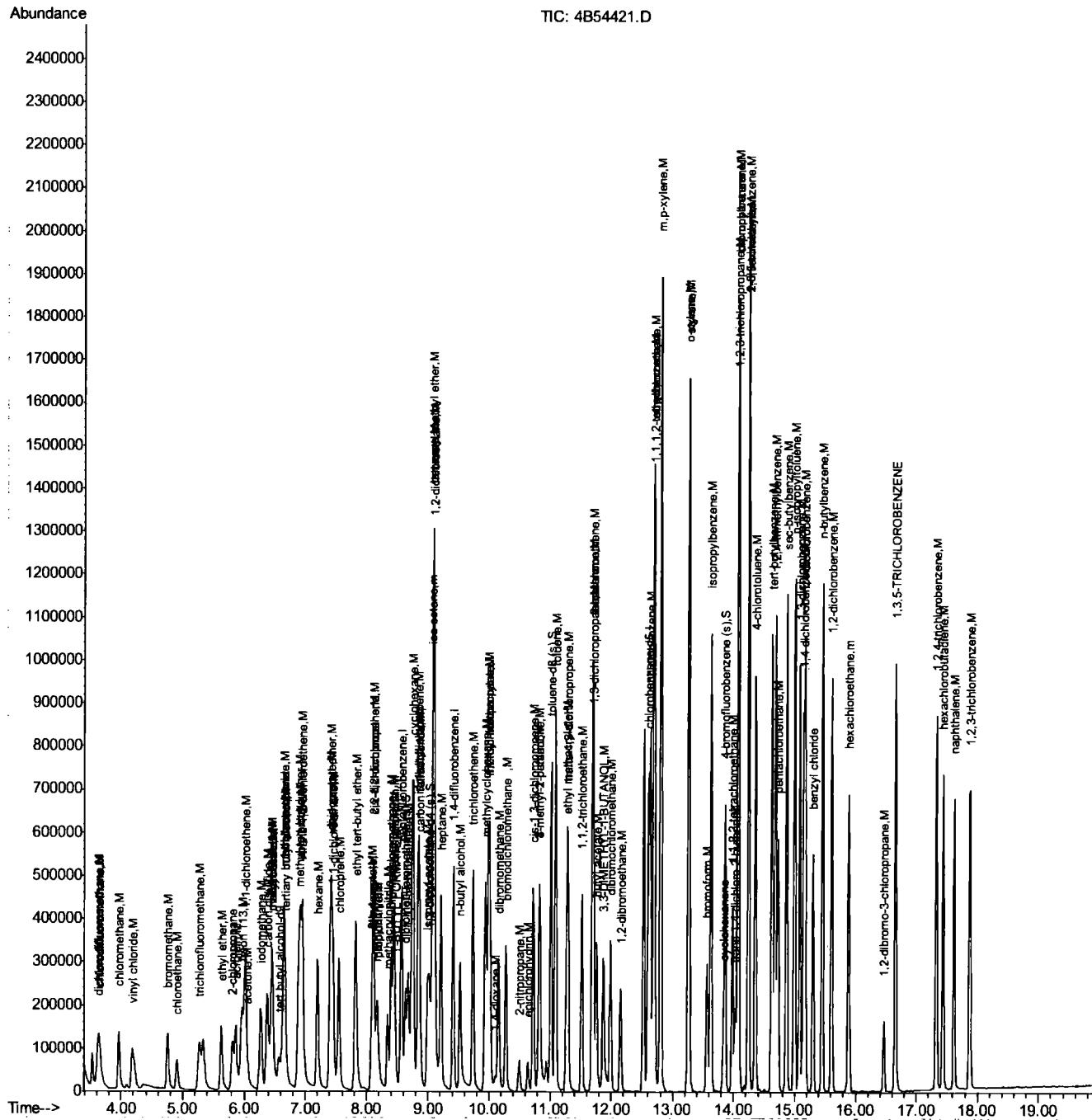
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
104) isopropylbenzene	13.63	105	837973	50.73	ug/L	99
106) cyclohexanone	13.83	55	67454	143.12	ug/L	96
107) bromobenzene	14.07	156	252617	52.14	ug/L	94
108) 1,1,2,2-tetrachloroethane	13.98	83	254860	52.69	ug/L	99
109) trans-1,4-dichloro-2-buten	14.02	53	39860	32.05	ug/L	97
110) 1,2,3-trichloropropane	14.06	110	60736	52.99	ug/L	99
111) n-propylbenzene	14.08	91	940003	50.35	ug/L	99
113) 2-chlorotoluene	14.24	126	214132	50.60	ug/L	93
114) 4-chlorotoluene	14.35	91	608802	50.43	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	707216	50.86	ug/L	99
116) tert-butylbenzene	14.62	119	640203	53.08	ug/L	98
117) pentachloroethane	14.72	167	149461	59.88	ug/L	99
118) 1,2,4-trimethylbenzene	14.68	105	703004	52.49	ug/L	100
119) sec-butylbenzene	14.87	105	953129	52.01	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	479023	54.31	ug/L	99
121) p-isopropyltoluene	15.00	119	815334	53.22	ug/L	100
122) 1,4-dichlorobenzene	15.17	146	470985	53.39	ug/L	99
123) benzyl chloride	15.30	91	448143	50.83	ug/L	98
124) 1,2-dichlorobenzene	15.61	146	467546	54.29	ug/L	99
126) n-butylbenzene	15.46	92	406730	53.53	ug/L	99
128) 1,2-dibromo-3-chloropropan	16.46	75	41838	53.89	ug/L	96
129) 1,3,5-TRICHLOROBENZENE	16.65	180	401126	58.40	ug/L	100
130) 1,2,4-trichlorobenzene	17.33	180	320514	58.26	ug/L	98
131) hexachlorobutadiene	17.44	225	196283	58.62	ug/L	99
132) naphthalene	17.62	128	574398	59.39	ug/L	100
133) 1,2,3-trichlorobenzene	17.88	180	269989	58.19	ug/L	99
134) hexachloroethane	15.89	201	167337	52.67	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54421.D
Acq On : 12 Aug 2015 4:55 pm
Operator : TOANP
Sample : jc1107-3msd
Misc : MS89470,V4B2296,w,,,1
ALS Vial : 17 Sample Multiplier: 1

Quant Time: Aug 12 17:34:44 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54459.D
 Acq On : 13 Aug 2015 11:45 am
 Operator : TOANP
 Sample : JC869-4MS
 Misc : MS89344, V4B2297, w, , , 1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 13 16:15:51 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	137687	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	412721	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	476155	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	427821	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	246362	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	150570	50.45	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 100.90%	
50) 1,2-dichloroethane-d4 (s)	9.01	65	155487	47.41	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 94.82%	
80) toluene-d8 (s)	11.02	98	552932	49.33	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 98.66%	
105) 4-bromofluorobenzene (s)	13.86	95	209256	48.65	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 97.30%	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) tertiary butyl alcohol	6.67	59	183278	550.29	ug/L	# 47
3) 1,4-dioxane	10.11	88	41734	1355.38	ug/L	92
7) chlorodifluoromethane	3.64	51	135136	32.38	ug/L	91
8) dichlorodifluoromethane	3.61	85	262049	65.56	ug/L	98
10) chloromethane	3.97	52	77549	50.73	ug/L	98
11) vinyl chloride	4.18	62	239658	54.56	ug/L	99
12) bromomethane	4.76	94	131315	52.13	ug/L	97
13) chloroethane	4.91	64	97826	56.64	ug/L	97
15) trichlorofluoromethane	5.28	101	287855	69.37	ug/L	99
19) ethyl ether	5.63	74	72543	46.13	ug/L	89
20) 2-chloropropane	5.81	39	29441	58.83	ug/L	90
21) acrolein	5.86	56	190735	362.20	ug/L	98
22) 1,1-dichloroethene	6.02	96	177639	54.51	ug/L	93
23) acetone	6.05	58	12921	47.13	ug/L	# 80
24) allyl chloride	6.46	76	104117	43.64	ug/L	# 68
25) acetonitrile	6.45	40	118532	451.62	ug/L	# 82
26) iodomethane	6.27	142	336270	54.65	ug/L	99
27) carbon disulfide	6.38	76	569086	59.51	ug/L	99
28) methylene chloride	6.64	84	175232	50.76	ug/L	90
29) methyl acetate	6.44	74	22891	49.27	ug/L	# 83
30) 1-chloropropane	6.66	42	284968	48.37	ug/L	# 95
31) methyl tert butyl ether	6.89	73	491494	56.52	ug/L	97
32) trans-1,2-dichloroethene	6.95	96	165879	51.75	ug/L	94
33) di-isopropyl ether	7.41	45	507107	44.47	ug/L	81
34) 2-butanone	8.08	72	17548	48.04	ug/L	# 90
35) 1,1-dichloroethane	7.45	63	294914	48.96	ug/L	98
36) chloroprene	7.55	53	227088	50.91	ug/L	97
37) acrylonitrile	6.92	53	268714	219.06	ug/L	99
38) vinyl acetate	7.43	86	23969	51.51	ug/L	82
39) ethyl tert-butyl ether	7.82	59	495070	48.87	ug/L	97
40) ethyl acetate	8.08	45	20488	41.32	ug/L	73
41) 2,2-dichloropropane	8.11	77	142835	56.09	ug/L	97
42) cis-1,2-dichloroethene	8.11	96	180575	51.86	ug/L	93

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54459.D
 Acq On : 13 Aug 2015 11:45 am
 Operator : TOANP
 Sample : JC869-4MS
 Misc : MS89344,V4B2297,w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 13 16:15:51 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.16	85	22868	48.36	ug/L	# 83
44) propionitrile	8.18	54	225830	456.44	ug/L	96
45) bromochloromethane	8.40	128	96370	53.93	ug/L	87
46) tetrahydrofuran	8.43	42	47851	40.25	ug/L	97
47) chloroform	8.44	85	190991	53.00	ug/L	98
48) T-BUTYL FORMATE	8.46	59	104446	45.34	ug/L	99
51) freon 113	5.97	151	115366	58.31	ug/L	89
52) methacrylonitrile	8.34	41	92060	42.13	ug/L	96
53) 1,1,1-trichloroethane	8.67	97	224337	58.87	ug/L	97
54) cyclohexane	8.74	84	235781	56.46	ug/L	91
55) iso-butyl alcohol	8.83	43	64841	476.85	ug/L	96
57) epichlorohydrin	10.63	57	72658	222.27	ug/L	98
58) n-butyl alcohol	9.53	56	218861	2361.10	ug/L	96
59) carbon tetrachloride	8.86	117	209958	55.11	ug/L	100
60) 1,1-dichloropropene	8.83	75	228712	54.03	ug/L	99
61) hexane	7.20	57	170260	52.95	ug/L	96
63) benzene	9.08	78	634983	50.96	ug/L	99
64) iso-octane	9.07	57	503338	50.29	ug/L	99
65) tert-amyl methyl ether	9.09	87	99530	50.24	ug/L	95
66) heptane	9.21	57	158642	71.66	ug/L	98
67) isopropyl acetate	8.99	61	61837	41.51	ug/L	98
68) 1,2-dichloroethane	9.10	62	213003	51.88	ug/L	95
69) trichloroethene	9.74	95	184757	55.09	ug/L	97
72) 2-nitropropane	10.50	41	45116	33.56	ug/L	# 1
73) 2-chloroethyl vinyl ether	10.50	63	23289	11.72	ug/L	96
74) methyl methacrylate	9.99	100	41578	52.03	ug/L	95
75) 1,2-dichloropropane	10.00	63	174541	48.76	ug/L	99
76) dibromomethane	10.16	93	109586	52.57	ug/L	96
77) methylcyclohexane	9.94	83	262192	60.80	ug/L	98
78) bromodichloromethane	10.28	83	241463	53.87	ug/L	100
79) cis-1,3-dichloropropene	10.72	75	291094	50.12	ug/L	99
81) 4-methyl-2-pentanone	10.82	58	63500	46.15	ug/L	96
82) toluene	11.09	92	422525	53.20	ug/L	100
83) 3-methyl-1-butanol	10.83	55	144751	942.29	ug/L	93
84) trans-1,3-dichloropropene	11.29	75	263178	51.74	ug/L	96
85) ethyl methacrylate	11.27	69	207396	49.29	ug/L	94
86) 1,1,2-trichloroethane	11.52	83	135665	49.07	ug/L	98
87) 2-hexanone	11.69	58	55327	44.04	ug/L	89
89) tetrachloroethene	11.69	164	204431	55.00	ug/L	98
90) 1,3-dichloropropane	11.71	76	249269	49.50	ug/L	96
91) butyl acetate	11.76	56	101836	44.34	ug/L	92
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	162916	483.38	ug/L	97
93) dibromochloromethane	11.99	129	206021	52.32	ug/L	99
94) 1,2-dibromoethane	12.15	107	173460	52.59	ug/L	100
96) chlorobenzene	12.63	112	492295	54.41	ug/L	97
97) 1,1,1,2-tetrachloroethane	12.70	131	182770	52.85	ug/L	100
98) ethylbenzene	12.68	91	822760	54.59	ug/L	98
99) m,p-xylene	12.80	106	632042	110.06	ug/L	99
100) o-xylene	13.25	106	320643	55.38	ug/L	97
101) styrene	13.27	104	531791	53.79	ug/L	96

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Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54459.D
 Acq On : 13 Aug 2015 11:45 am
 Operator : TOANP
 Sample : JC869-4MS
 Misc : MS89344,V4B2297,w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 13 16:15:51 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

743 7

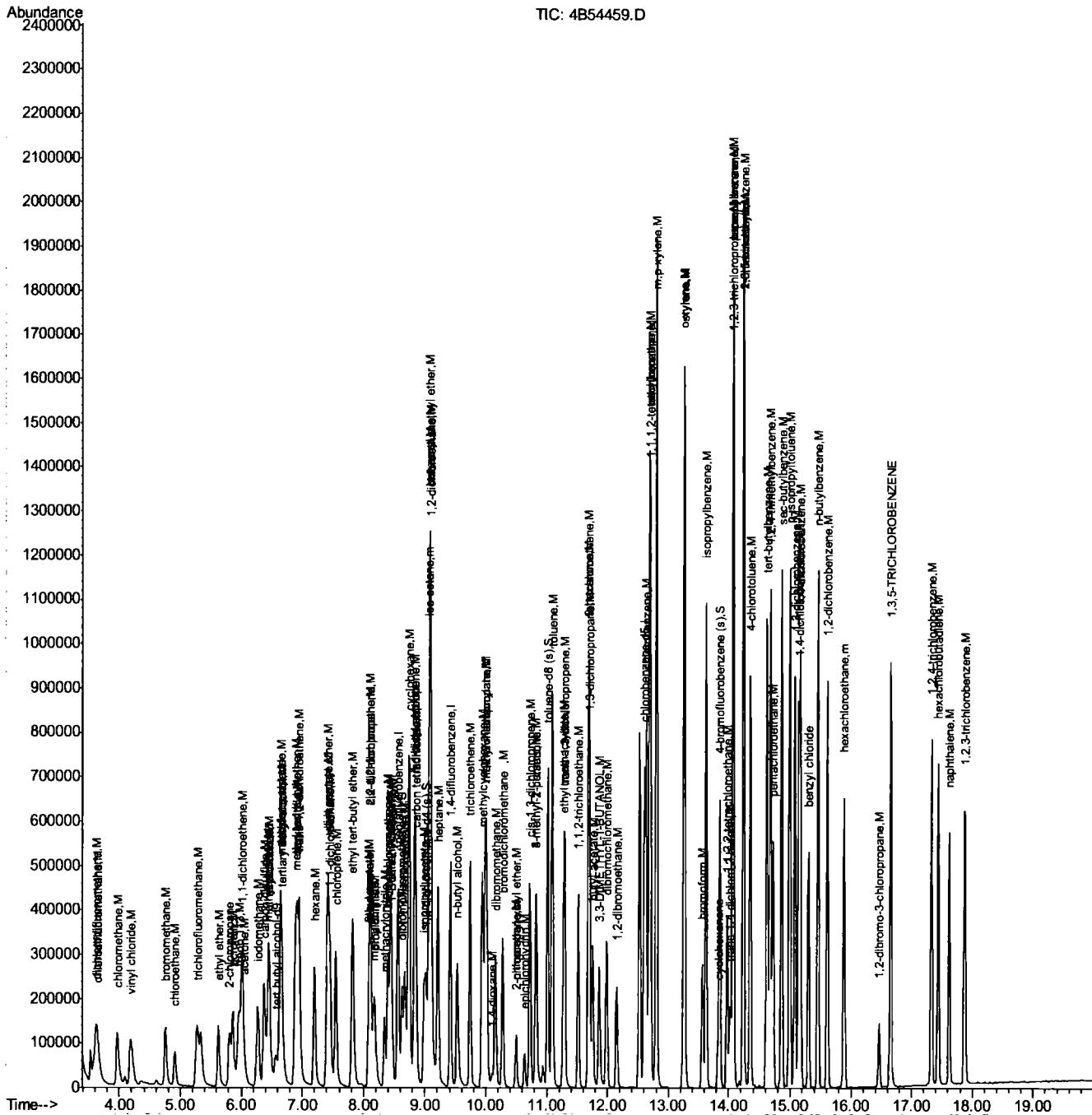
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
102) bromoform	13.57	173	157134	51.35	ug/L	98
104) isopropylbenzene	13.63	105	843758	53.95	ug/L	99
106) cyclohexanone	13.82	55	56354	126.29	ug/L	96
107) bromobenzene	14.07	156	246694	53.78	ug/L	92
108) 1,1,2,2-tetrachloroethane	13.98	83	222670	48.62	ug/L	98
109) trans-1,4-dichloro-2-butene	14.02	53	38683	32.85	ug/L	93
110) 1,2,3-trichloropropane	14.06	110	54928	50.62	ug/L	96
111) n-propylbenzene	14.08	91	942196	53.30	ug/L	98
113) 2-chlorotoluene	14.24	126	210757	52.60	ug/L	95
114) 4-chlorotoluene	14.35	91	596143	52.16	ug/L	100
115) 1,3,5-trimethylbenzene	14.24	105	697302	52.97	ug/L	99
116) tert-butylbenzene	14.62	119	632393	55.38	ug/L	99
117) pentachloroethane	14.72	167	141496	59.87	ug/L	97
118) 1,2,4-trimethylbenzene	14.68	105	682393	53.81	ug/L	99
119) sec-butylbenzene	14.87	105	953850	54.97	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	455976	54.60	ug/L	99
121) p-isopropyltoluene	15.00	119	807374	55.66	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	446086	53.41	ug/L	99
123) benzyl chloride	15.30	91	429949	51.50	ug/L	98
124) 1,2-dichlorobenzene	15.61	146	441998	54.21	ug/L	98
126) n-butylbenzene	15.46	92	395869	55.02	ug/L	99
128) 1,2-dibromo-3-chloropropan	16.46	75	35726	48.60	ug/L	93
129) 1,3,5-TRICHLOROBENZENE	16.65	180	373244	57.40	ug/L	99
130) 1,2,4-trichlorobenzene	17.33	180	282580	54.25	ug/L	98
131) hexachlorobutadiene	17.44	225	190359	60.04	ug/L	99
132) naphthalene	17.62	128	473209	51.68	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	236221	53.77	ug/L	99
134) hexachloroethane	15.89	201	157289	52.29	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54459.D
 Acq On : 13 Aug 2015 11:45 am
 Operator : TOANP
 Sample : JC869-4MS
 Misc : MS89344, V4B2297, w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 13 16:15:51 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54458.D
 Acq On : 13 Aug 2015 11:17 am
 Operator : TOANP
 Sample : JC869-3DUP
 Misc : MS89344,V4B2297,w,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 13 16:15:38 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	129185	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	418519	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	473275	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	402037	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	185731	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	150429	49.71	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery =	99.42%		
50) 1,2-dichloroethane-d4 (s)	9.02	65	152325	45.80	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery =	91.60%		
80) toluene-d8 (s)	11.02	98	539166	48.39	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery =	96.78%		
105) 4-bromofluorobenzene (s)	13.86	95	177414	54.71	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	109.42%		

Target Compounds

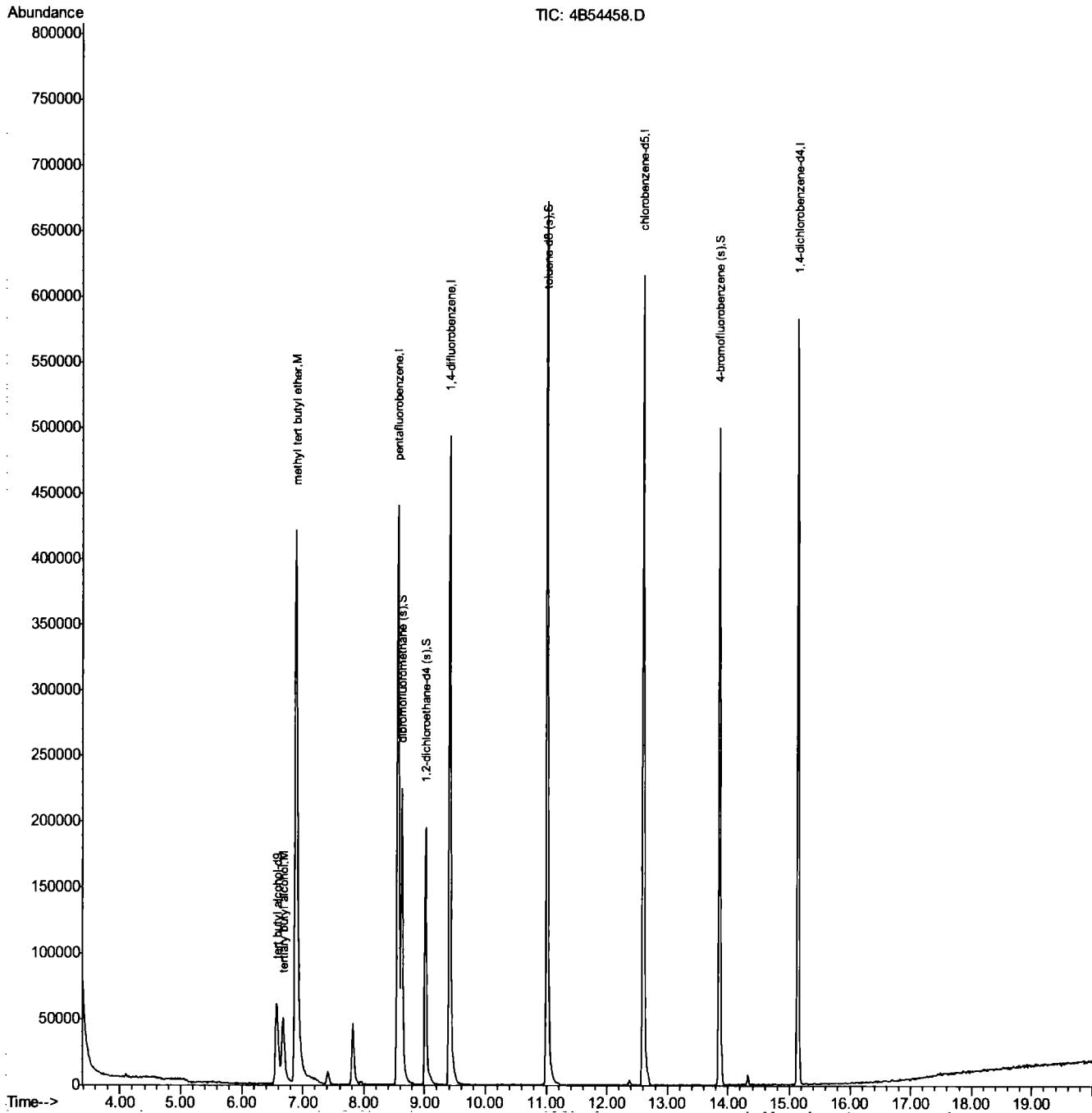
					Qvalue
2) tertiary butyl alcohol	6.68	59	99984	319.96	ug/L # 20
31) methyl tert butyl ether	6.89	73	674325	76.48	ug/L 98

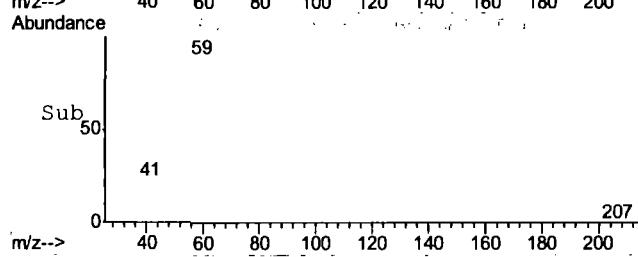
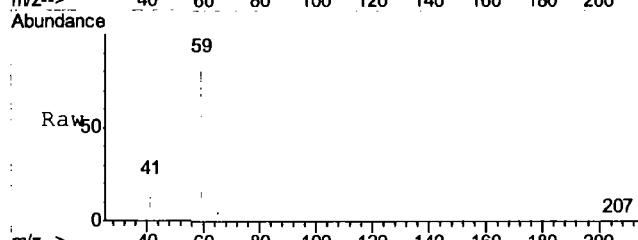
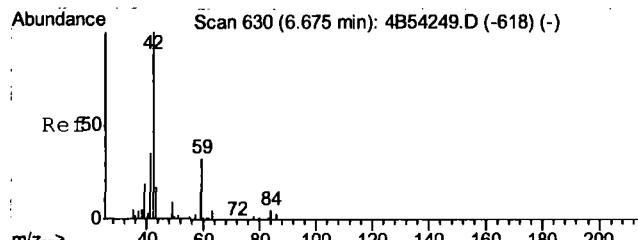
(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54458.D
Acq On : 13 Aug 2015 11:17 am
Operator : TOANP
Sample : JC869-3DUP
Misc : MS89344, V4B2297, w,,,1
ALS Vial : 5 Sample Multiplier: 1

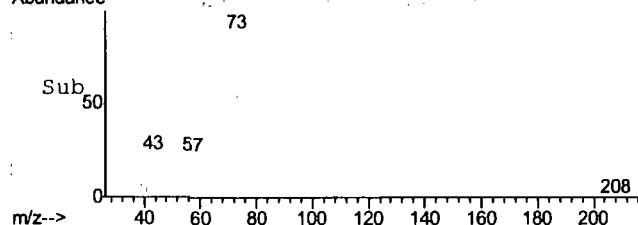
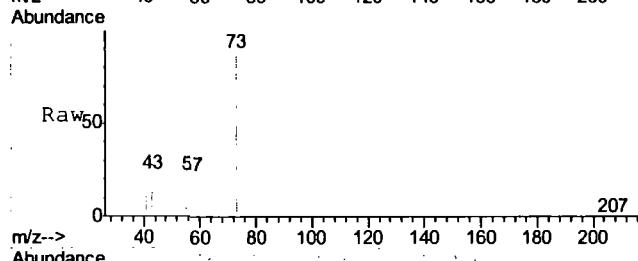
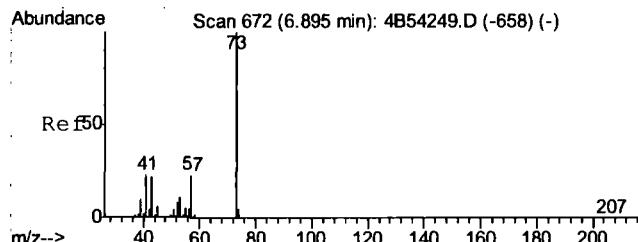
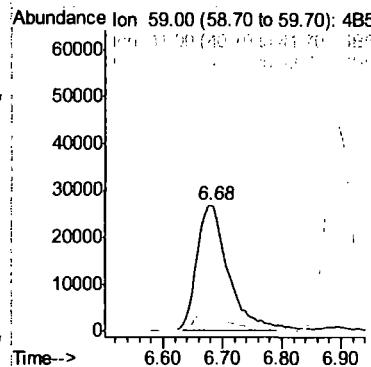
Quant Time: Aug 13 16:15:38 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration





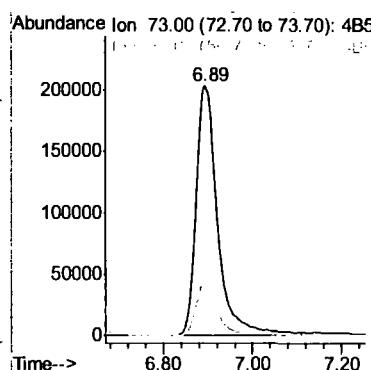
#2
tertiary butyl alcohol
Concen: 319.96 ug/L
RT: 6.68 min Scan# 631
Delta R.T. 0.01 min
Lab File: 4B54458.D
Acq: 13 Aug 2015 11:17 am

Tgt Ion: 59 Resp: 99984
Ion Ratio Lower Upper
59 100
41 22.0 92.9 152.9#
43 12.9 25.2 85.2#



#31
methyl tert butyl ether
Concen: 76.48 ug/L
RT: 6.89 min Scan# 672
Delta R.T. -0.00 min
Lab File: 4B54458.D
Acq: 13 Aug 2015 11:17 am

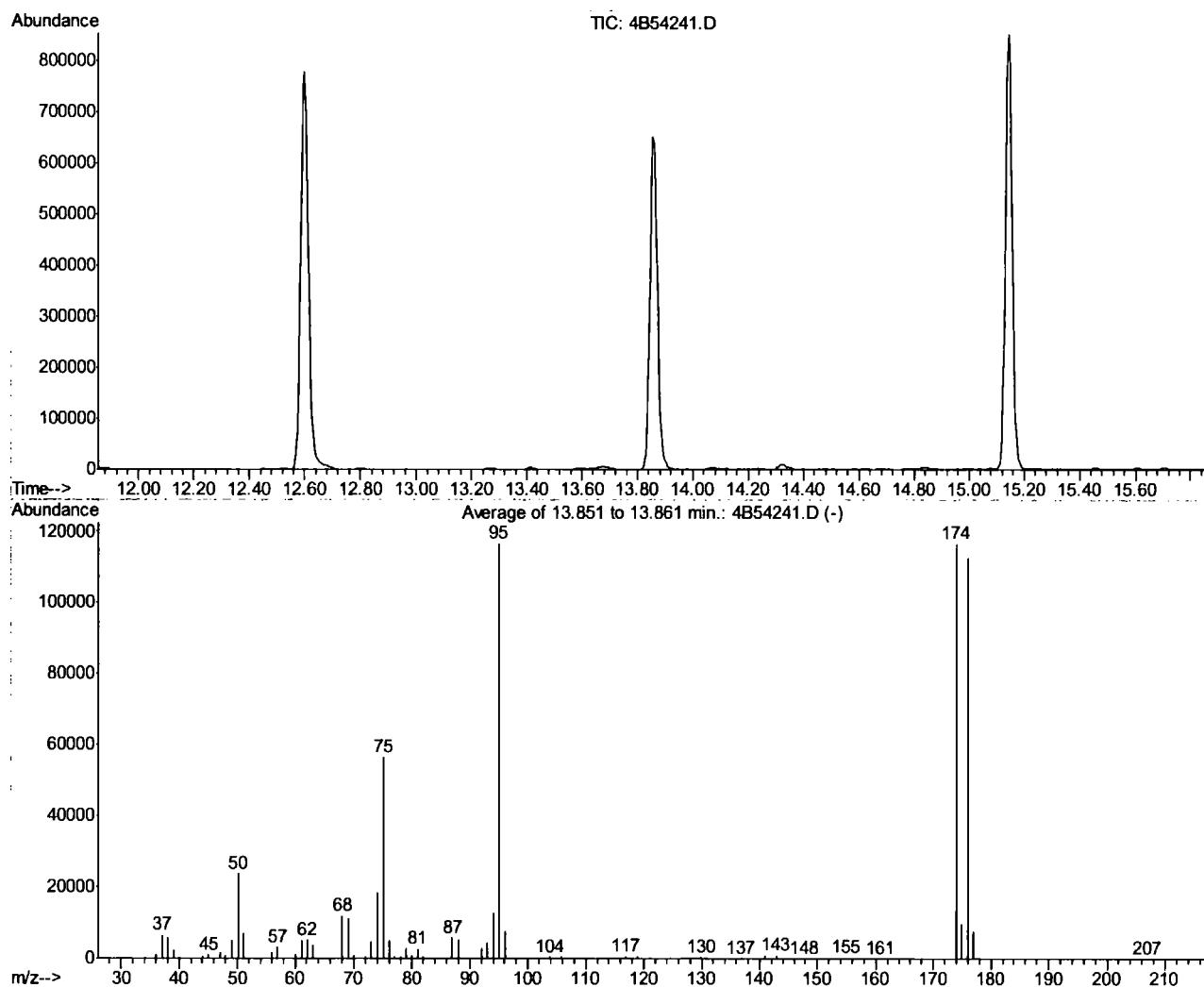
Tgt Ion: 73 Resp: 674325
Ion Ratio Lower Upper
73 100
57 22.0 0.0 68.7



SW-846 Method 8260

Data File : C:\MSDCHEM\1\DATA\4B54241.D Vial: 1
 Acq On : 5 Aug 2015 9:05 am Operator: TOANP
 Sample : bfb Inst : MS4B
 Misc : MS88663,V4B2289,w,,,1 Multiplr: 1.00
 MS Integration Params: RTEINT.P

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um



AutoFind: Scans 2002, 2003, 2004; Background Corrected with Scan 1993

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	20.4	23821	PASS
75	95	30	60	48.5	56474	PASS
95	95	100	100	100.0	116525	PASS
96	95	5	9	6.8	7868	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	100.1	116618	PASS
175	174	5	9	8.4	9749	PASS
176	174	95	101	96.6	112621	PASS
177	176	5	9	7.0	7858	PASS

Average of 13.851 to 13.861 min.: 4B54241.D

fbf

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.10	1315	51.10	7161	64.00	116	77.00	751
37.10	6644	52.10	338	64.20	181	78.00	544
38.10	5932	55.10	458	67.10	142	78.95	2868
39.10	2254	56.05	1766	68.00	11950	79.95	960
40.00	273	57.05	3179	69.10	11507	81.00	2838
44.05	634	57.90	72	69.95	843	81.90	627
45.05	1108	58.10	60	72.05	612	86.00	110
47.05	1902	60.00	1094	73.00	4764	87.00	5892
48.00	795	61.10	5023	74.10	18488	88.00	5478
49.10	5206	62.10	5332	75.10	56474	90.95	382
50.10	23821	63.10	3765	76.05	5002	92.05	2969

Average of 13.851 to 13.861 min.: 4B54241.D

fbf

Modified:subtracted

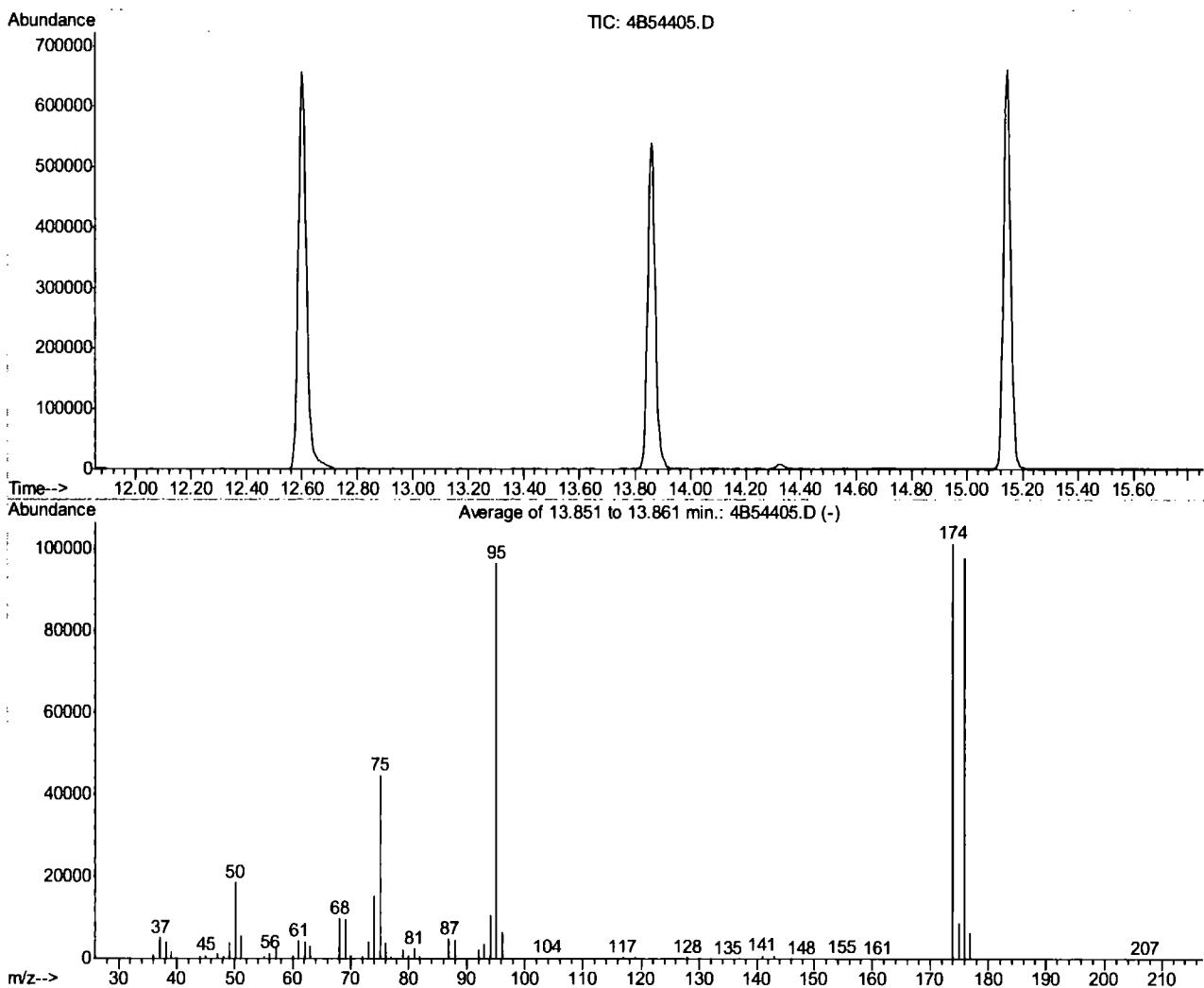
m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
93.05	4444	117.95	344	142.95	1027	173.95	116618
94.05	12802	118.95	544	145.70	62	175.00	9749
95.00	116525	127.90	458	147.00	55	175.95	112621
96.05	7868	128.95	162	147.70	66	176.95	7858
96.95	221	129.95	495	147.90	214	177.90	224
103.90	528	130.90	194	149.00	57	207.00	54
104.90	60	134.95	113	149.90	54		
105.10	58	136.85	222	155.05	313		
105.90	506	139.80	52	156.95	219		
115.95	399	141.00	861	160.90	55		
116.90	674	142.00	99	171.30	51		

SW-846 Method 8260

Data File : C:\MSDCHEM\1\DATA\4B54405.D
 Acq On : 12 Aug 2015 9:13 am
 Sample : BFB
 Misc : MS89342, V4B2296, w,,,1
 MS Integration Params: RTEINT.P

Vial: 1
 Operator: TOANP
 Inst : MS4B
 Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um



AutoFind: Scans 2002, 2003, 2004; Background Corrected with Scan 1993

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	19.3	18701	PASS
75	95	30	60	46.4	44842	PASS
95	95	100	100	100.0	96672	PASS
96	95	5	9	6.7	6517	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	105.0	101530	PASS
175	174	5	9	8.6	8723	PASS
176	174	95	101	96.6	98101	PASS
177	176	5	9	6.7	6561	PASS

Average of 13.851 to 13.861 min.: 4B54405.D

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.05	998	49.10	3996	64.05	268	77.90	306
37.10	5202	50.10	18701	67.05	285	78.10	195
38.10	4225	51.10	5773	68.00	9862	78.95	2347
39.10	1750	52.15	263	69.05	9506	80.00	683
40.00	215	55.05	426	70.05	742	80.95	2506
42.05	125	56.05	1391	72.05	648	81.95	481
43.20	54	57.10	2507	73.05	4064	86.00	64
44.00	489	60.05	781	74.05	15256	87.00	4883
45.05	803	61.05	4351	75.10	44842	88.00	4597
47.00	1356	62.10	4225	76.05	3933	91.00	320
48.10	633	63.10	3051	76.95	629	92.00	2436

Average of 13.851 to 13.861 min.: 4B54405.D

BFB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
93.00	3623	117.90	242	142.95	876	174.00	101530
94.10	10563	118.95	495	145.00	54	175.00	8723
95.05	96672	127.95	413	145.90	142	175.95	98101
96.05	6517	128.90	135	147.00	54	176.95	6561
97.10	210	129.95	384	147.95	228	177.90	255
103.90	374	130.90	52	149.95	114	207.15	149
104.90	130	135.00	120	154.95	287		
106.00	301	136.90	145	156.95	205		
106.90	52	140.95	882	159.00	61		
115.95	331	142.00	57	160.95	103		
116.90	513	142.30	66	172.25	238		

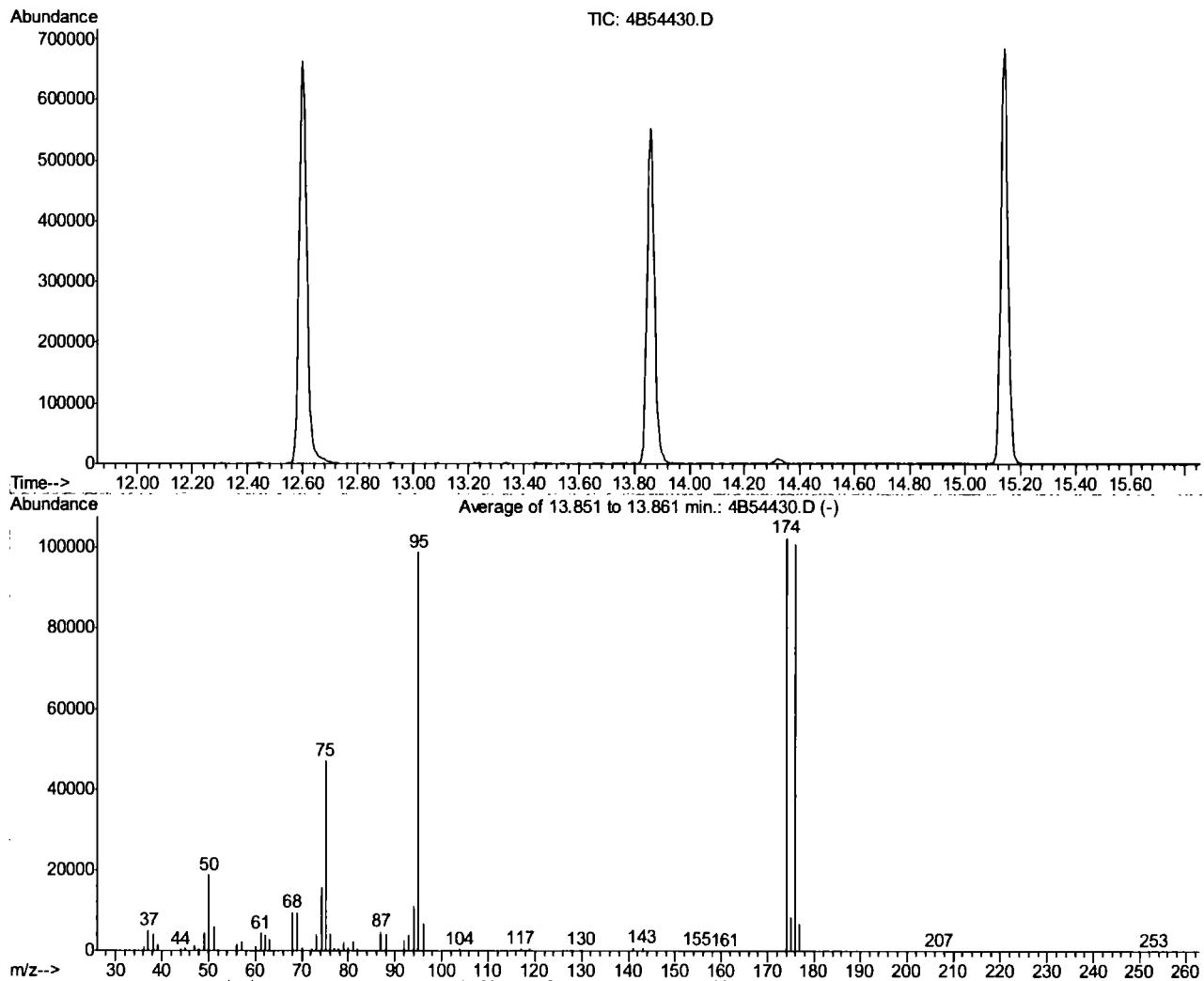
7.62 7

SW-846 Method 8260

Data File : C:\MSDCHEM\1\DATA\4B54430.D
 Acq On : 12 Aug 2015 9:06 pm
 Sample : bfb
 Misc : MS89499,V4B2297,w,,,1
 MS Integration Params: RTEINT.P

Vial: 26
 Operator: TOANP
 Inst : MS4B
 Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um



AutoFind: Scans 2002, 2003, 2004; Background Corrected with Scan 1993

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
50	95	15	40	19.0	18824	PASS
75	95	30	60	47.6	47162	PASS
95	95	100	100	100.0	98989	PASS
96	95	5	9	6.8	6769	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	103.6	102565	PASS
175	174	5	9	8.2	8410	PASS
176	174	95	101	98.4	100890	PASS
177	176	5	9	6.7	6717	PASS

Average of 13.851 to 13.861 min.: 4B54430.D

bfB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.10	1012	51.10	6094	67.10	98	78.95	2224
37.10	5056	52.05	251	68.05	9518	79.95	756
38.10	4108	55.10	253	69.00	9514	80.95	2371
39.10	1608	56.05	1467	70.05	735	81.95	486
40.00	61	57.05	2415	72.15	473	85.95	130
44.05	475	58.10	57	72.95	3843	87.00	4857
45.05	880	60.05	967	74.10	15718	88.00	4310
47.05	1324	61.10	4403	75.10	47162	91.00	263
48.05	634	62.05	4068	76.10	4189	91.20	107
49.10	4499	63.05	3017	77.05	712	92.05	2676
50.10	18824	64.05	301	77.95	480	93.00	3941

Average of 13.851 to 13.861 min.: 4B54430.D

bfB

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
94.05	11051	117.80	321	146.90	57	177.90	219
95.00	98989	118.95	435	147.95	258	206.95	116
96.05	6769	127.85	324	154.90	306	253.00	157
97.05	140	129.00	131	156.95	134		
103.95	430	129.90	377	160.90	50		
104.95	122	130.90	70	171.75	118		
105.95	388	134.90	188	172.20	98		
106.90	52	136.95	107	174.00	102565		
115.80	64	140.95	867	175.00	8410		
115.95	168	141.90	75	176.00	100890		
117.00	574	142.95	884	176.90	6717		

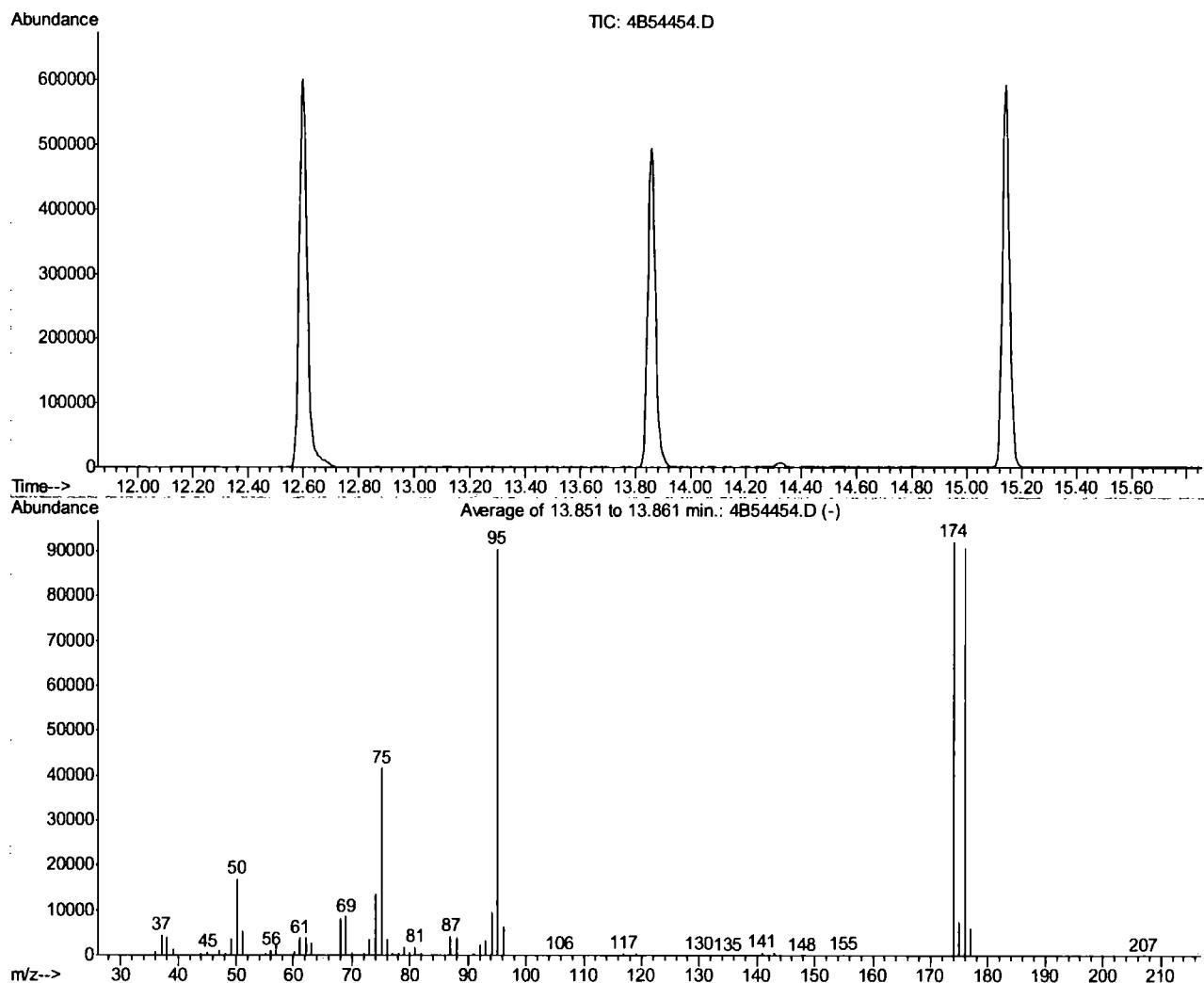
763
7

SW-846 Method 8260

Data File : C:\MSDCHEM\1\DATA\4B54454.D
 Acq On : 13 Aug 2015 9:00 am
 Sample : bfb
 Misc : MS89470, V4B2298, w, , , 1
 MS Integration Params: RTEINT.P

Vial: 1
 Operator: TOANP
 Inst : MS4B
 Multiplr: 1.00

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um



AutoFind: Scans 2002, 2003, 2004; Background Corrected with Scan 1993

Target	Rel. to	Lower	Upper	Rel.	Raw	Result
Mass	Mass	Limit%	Limit%	Abn%	Abn	Pass/Fail
50	95	15	40	18.8	17050	PASS
75	95	30	60	46.1	41786	PASS
95	95	100	100	100.0	90570	PASS
96	95	5	9	7.0	6316	PASS
173	174	0.00	2	0.0	0	PASS
174	95	50	120	101.9	92250	PASS
175	174	5	9	8.2	7549	PASS
176	174	95	101	98.3	90642	PASS
177	176	5	9	6.8	6125	PASS

Average of 13.851 to 13.861 min.: 4B54454.D

bf
bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
36.10	862	50.10	17050	67.10	110	79.00	1922
37.10	4563	51.10	5337	68.05	8334	79.95	624
38.10	3942	52.05	262	69.00	8693	80.95	1965
39.10	1502	55.05	374	70.05	686	82.00	477
40.00	33	56.10	1273	72.05	478	87.00	4259
44.00	528	57.05	2298	73.00	3548	88.00	3952
45.10	803	60.05	937	74.10	13782	91.00	371
47.05	1177	61.05	3968	75.10	41786	92.00	2306
47.80	136	62.10	3929	76.05	3601	93.05	3393
48.15	526	63.10	2811	77.00	574	94.10	9760
49.10	3685	64.10	235	78.05	426	95.00	90570

Average of 13.851 to 13.861 min.: 4B54454.D

bf
bfb

Modified:subtracted

m/z	abund.	m/z	abund.	m/z	abund.	m/z	abund.
96.05	6316	128.95	205	154.95	247		
97.00	230	129.95	320	156.90	193		
103.95	358	134.95	197	158.80	50		
104.95	126	136.90	120	172.10	74		
105.90	361	140.95	759	174.00	92250		
107.00	59	143.00	706	175.00	7549		
115.95	314	145.90	50	176.00	90642		
116.95	502	147.90	238	176.95	6125		
117.95	302	150.00	56	177.90	127		
119.00	501	152.90	50	178.10	64		
127.95	301	153.90	62	207.05	170		

764

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54242.D
 Acq On : 5 Aug 2015 9:39 am
 Operator : TOANP
 Sample : ic2289-0.2
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 07 10:38:36 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:57:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	189317	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	455240	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	533050	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	471164	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	250843	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	165428	50.74	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	101.48%
50) 1,2-dichloroethane-d4 (s)	9.02	65	183485	50.72	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	101.44%
80) toluene-d8 (s)	11.02	98	626021	49.88	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	99.76%
105) 4-bromofluorobenzene (s)	13.86	95	227429	51.33	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	102.66%

Target Compounds

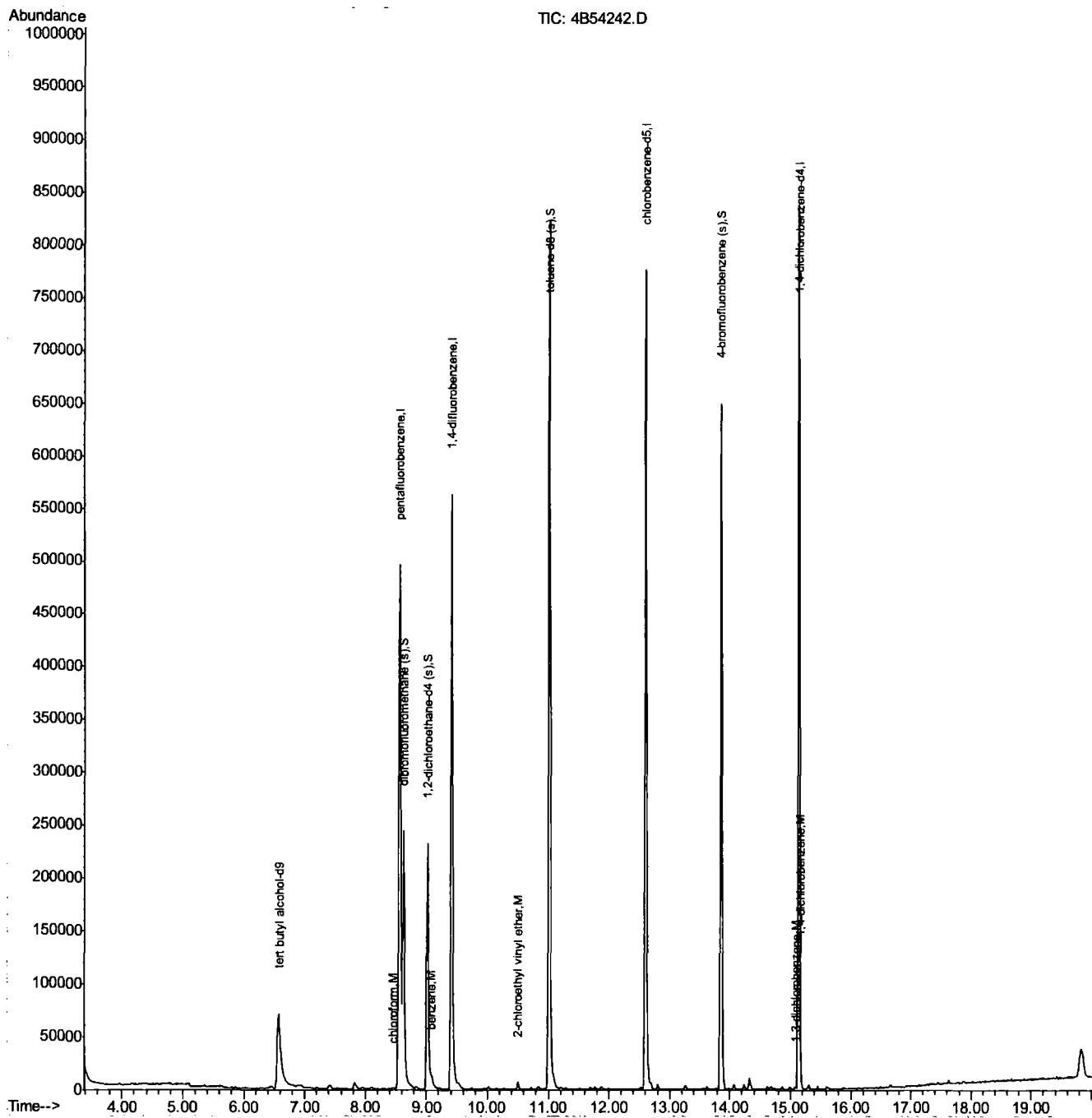
					Qvalue
47) chloroform	8.44	85	732	0.18	ug/L # 67
63) benzene	9.08	78	2867	0.20	ug/L 92
73) 2-chloroethyl vinyl ether	10.50	63	3836	1.76	ug/L 96
120) 1,3-dichlorobenzene	15.08	146	1473	0.17	ug/L 96
122) 1,4-dichlorobenzene	15.17	146	1583	0.18	ug/L 79

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54242.D
 Acq On : 5 Aug 2015 9:39 am
 Operator : TOANP
 Sample : ic2289-0.2
 Misc : MS88663, V4B2289, w,,,,,1
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 07 10:38:36 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:57:32 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Manual Integrations
APPROVED
(compounds with "m" flag)
Jessica Reitan-Chu
08/10/15 09:30

Data Path : C:\MSDChem\1\DATA\
Data File : 4B54243.D
Acq On : 5 Aug 2015 10:07 am
Operator : TOANP
Sample : ic2289-0.5
Misc : MS88663,V4B2289,w,,,1
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 06 08:44:03 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 13:57:32 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) tert butyl alcohol-d9	6.58	65	161627	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	464201	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	539990	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	476569	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	253504	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	163660	49.23	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 98.46%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	185165	50.19	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 100.38%	
80) toluene-d8 (s)	11.02	98	634463	49.90	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 99.80%	
105) 4-bromofluorobenzene (s)	13.86	95	225253	50.30	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.60%	

Target Compounds

				Qvalue	
10) chloromethane	3.96	52	875	0.51	ug/L # 76
11) vinyl chloride	4.16	62	2163	0.47	ug/L 84
12) bromomethane	4.76	94	1598	0.49	ug/L 85
13) chloroethane	4.92	64	979	0.50	ug/L # 44
19) ethyl ether	5.63	74	735	0.42	ug/L # 65
21) acrolein	5.87	56	2685	4.76	ug/L 93
22) 1,1-dichloroethene	6.03	96	1501	0.40	ug/L # 66
26) iodomethane	6.28	142	2790	0.41	ug/L 72
27) carbon disulfide	6.38	76	4291	0.40	ug/L 71
28) methylene chloride	6.65	84	1664	0.43	ug/L # 59
31) methyl tert butyl ether	6.90	73	4143	0.45	ug/L 93
32) trans-1,2-dichloroethene	6.95	96	1720	0.45	ug/L # 66
33) di-isopropyl ether	7.41	45	5508	0.39	ug/L 90
39) ethyl tert-butyl ether	7.82	59	4980	0.39	ug/L 96
44) propionitrile	8.20	54	2348m	4.29	ug/L
47) chloroform	8.44	85	1690	0.41	ug/L 95
52) methacrylonitrile	8.34	41	1130	0.46	ug/L 87
57) epichlorohydrin	10.64	57	755	1.91	ug/L 66
58) n-butyl alcohol	9.54	56	2498	18.71	ug/L 95
63) benzene	9.08	78	6286	0.43	ug/L 87
68) 1,2-dichloroethane	9.10	62	1829	0.42	ug/L 74
72) 2-nitropropane	10.50	41	644	0.41	ug/L # 49
73) 2-chloroethyl vinyl ether	10.50	63	5053	2.28	ug/L 94
76) dibromomethane	10.16	93	943	0.40	ug/L 90
78) bromodichloromethane	10.28	83	1925	0.41	ug/L 81
84) trans-1,3-dichloropropene	11.30	75	2239	0.42	ug/L 90
90) 1,3-dichloropropane	11.71	76	2158	0.41	ug/L 98
92) 3,3-DIMETHYL-1-BUTANOL	11.88	57	1664	3.49	ug/L 86
94) 1,2-dibromoethane	12.15	107	1474	0.44	ug/L 90
96) chlorobenzene	12.64	112	3985	0.41	ug/L 89
98) ethylbenzene	12.69	91	6598	0.41	ug/L 88
99) m,p-xylene	12.80	106	4961	0.83	ug/L 94
100) o-xylene	13.25	106	2401	0.40	ug/L # 77

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54243.D
 Acq On : 5 Aug 2015 10:07 am
 Operator : TOANP
 Sample : ic2289-0.5
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 06 08:44:03 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:57:32 2015
 Response via : Initial Calibration

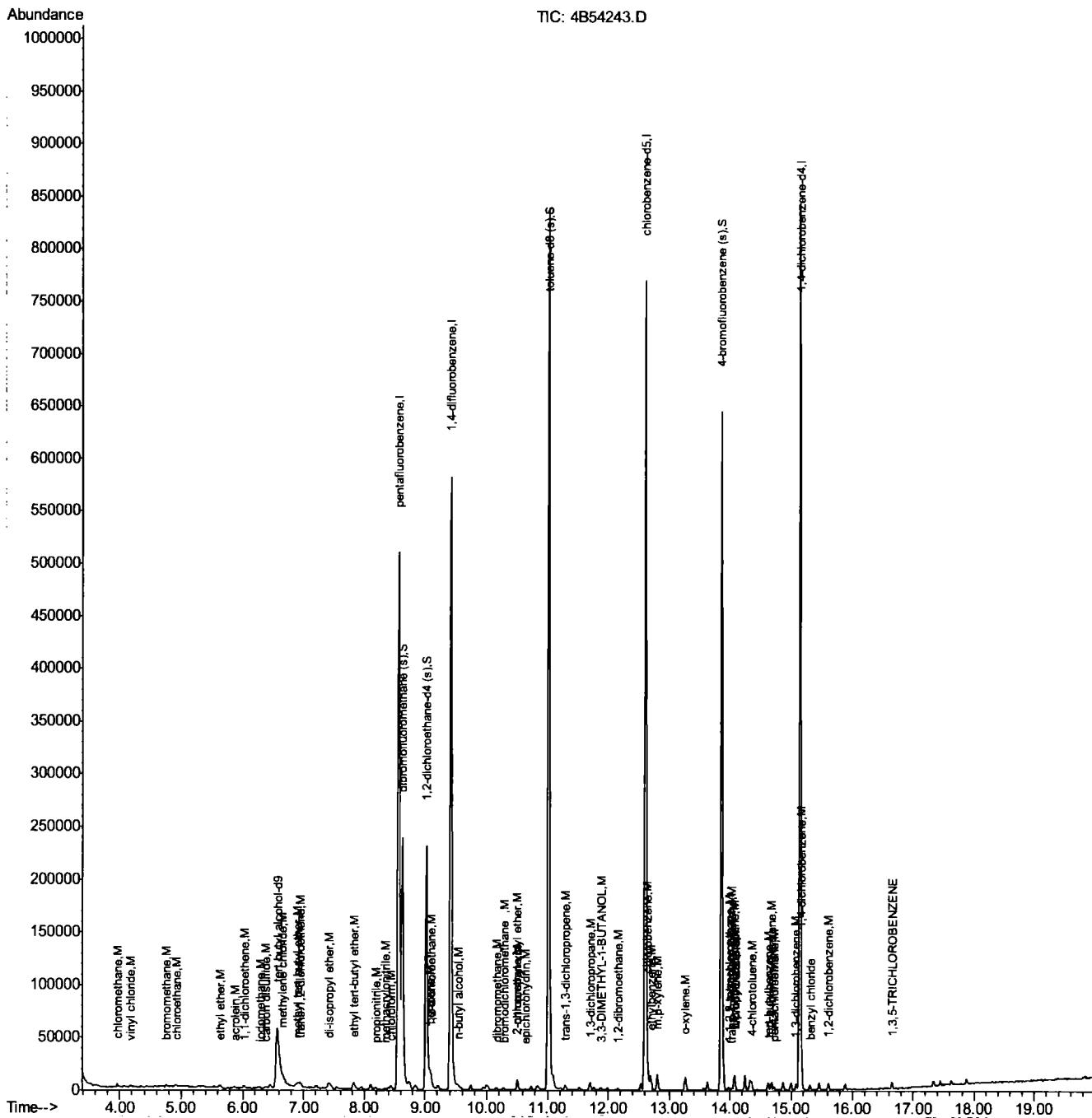
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
107) bromobenzene	14.07	156	1937	0.42	ug/L	94
108) 1,1,2,2-tetrachloroethane	13.98	83	2013	0.45	ug/L	94
109) trans-1,4-dichloro-2-buten	14.03	53	492	0.41	ug/L	# 64
110) 1,2,3-trichloropropane	14.07	110	455	0.41	ug/L	88
111) n-propylbenzene	14.08	91	7776	0.44	ug/L	99
114) 4-chlorotoluene	14.35	91	4846	0.42	ug/L	92
116) tert-butylbenzene	14.62	119	5430	0.48	ug/L	94
117) pentachloroethane	14.72	167	1029	0.42	ug/L	# 84
118) 1,2,4-trimethylbenzene	14.68	105	5277	0.42	ug/L	90
120) 1,3-dichlorobenzene	15.08	146	3591	0.42	ug/L	89
122) 1,4-dichlorobenzene	15.17	146	3728	0.43	ug/L	92
123) benzyl chloride	15.31	91	4240	0.42	ug/L	# 91
124) 1,2-dichlorobenzene	15.61	146	3345	0.41	ug/L	95
129) 1,3,5-TRICHLOROBENZENE	16.65	180	2681	0.42	ug/L	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
Data File : 4B54243.D
Acq On : 5 Aug 2015 10:07 am
Operator : TOANP
Sample : ic2289-0.5
Misc : MS88663,V4B2289,w,,,1
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 06 08:44:03 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 13:57:32 2015
Response via : Initial Calibration



Manual Integration Approval Summary

Page 1 of 1

Sample Number: V4B2289-IC2289 **Method:** SW846 8260C
Lab FileID: 4B54243.D **Analyst approved:** 08/06/15 09:13 Dong, Mei
Injection Time: 08/05/15 10:07 **Supervisor approved:** 08/10/15 09:30 Jessica Reitan-Chu

Parameter	CAS	Sig#	R.T. (min.)	Reason
Propionitrile	107-12-0		8.20	Poor instrument integration

7.7.2.1
7

Quantitation Report (Qedit)

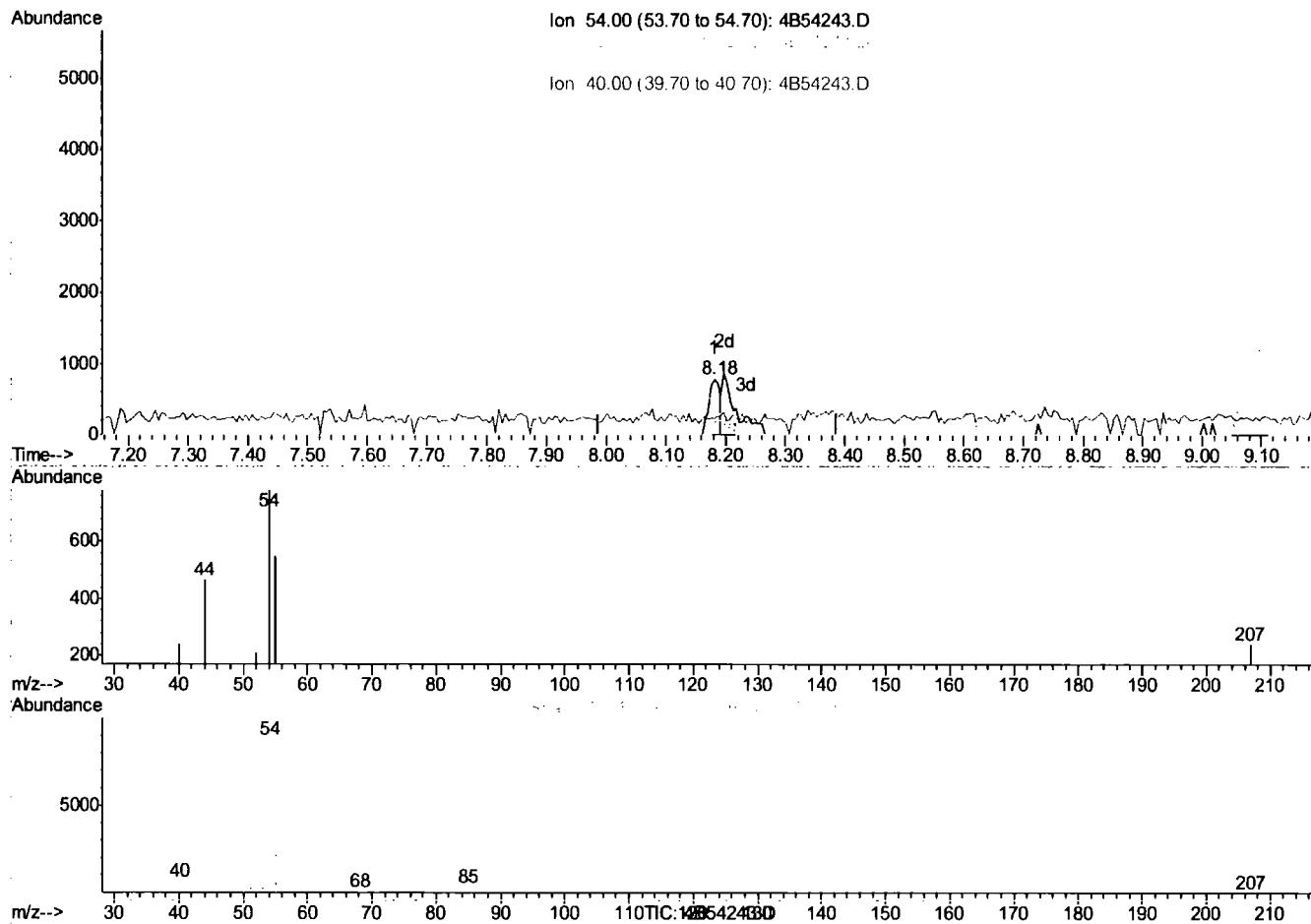
Data File : C:\MSDCHEM\1\DATA\4B54243.D
 Acq On : 5 Aug 2015 10:07 am
 Sample : ic2289-0.5
 Misc : MS88663,V4B2289,w,,,1
 MS Integration Params: RTEINT.P

Vial: 3
 Operator: TOANP
 Inst : MS4B
 Multiplr: 1.00

Quant Time: Aug 05 13:59:41 2015

Results File: M4B2289.RES

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 Last Update : Wed Aug 05 13:57:32 2015
 Response via : Multiple Level Calibration



(44) propionitrile (M)

8.18min 1.85ug/L

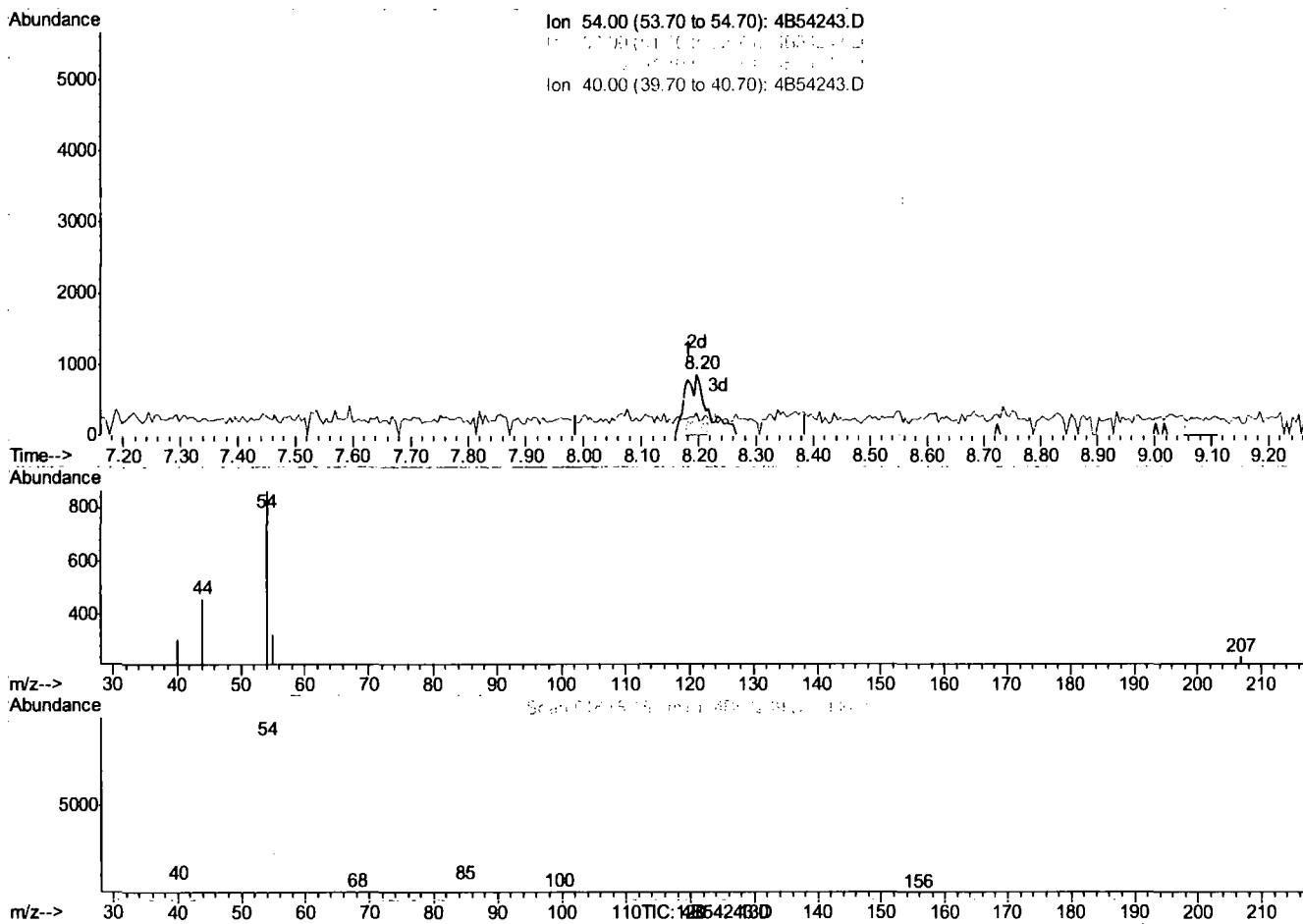
response 1014

Ion	Exp%	Act%
54.00	100	100
52.00	19.10	26.63
55.00	34.20	34.57
40.00	6.10	10.76

Quantitation Report (Qedit)
 Data File : C:\MSDChem\1\DATA\4B54243.D Vial: 3
 Acq On : 5 Aug 2015 10:07 am Operator: TOANP
 Sample : ic2289-0.5 Inst : MS4B
 Misc : MS88663,V4B2289,w,,,1 Multiplr: 1.00
 MS Integration Params: RTEINT.P

Quant Time: Aug 06 08:44:03 2015 Results File: M4B2289.RES

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 Last Update : Thu Aug 06 08:53:28 2015
 Response via : Multiple Level Calibration



(44) propionitrile (M)

8.20min 4.29ug/L m

response 2348

Ion	Exp%	Act%
54.00	100	100
52.00	19.10	24.88
55.00	34.20	37.85
40.00	6.10	35.88

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54244.D
 Acq On : 5 Aug 2015 10:35 am
 Operator : TOANF
 Sample : ic2289-1
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Aug 10 08:18:11 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:57:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	168106	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	456387	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	528116	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	470067	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	254497	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	163277	49.96	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	99.92%
50) 1,2-dichloroethane-d4 (s)	9.02	65	181441	50.02	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	100.04%
80) toluene-d8 (s)	11.02	98	621889	50.01	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	100.02%
105) 4-bromofluorobenzene (s)	13.86	95	223099	49.63	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.26%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.67	59	2271	5.49	ug/L # 72
7) chlorodifluoromethane	3.64	51	5246	1.20	ug/L 97
10) chloromethane	3.95	52	1654	0.97	ug/L # 61
11) vinyl chloride	4.16	62	4538	1.00	ug/L 85
12) bromomethane	4.76	94	2599	0.81	ug/L 92
13) chloroethane	4.92	64	1722	0.90	ug/L 86
15) trichlorofluoromethane	5.28	101	3571	0.86	ug/L # 62
19) ethyl ether	5.63	74	1897	1.10	ug/L 88
21) acrolein	5.87	56	5534	9.97	ug/L 80
22) 1,1-dichloroethene	6.01	96	4173	1.13	ug/L 86
24) allyl chloride	6.45	76	2918	1.11	ug/L # 33
26) iodomethane	6.28	142	7104	1.05	ug/L 89
27) carbon disulfide	6.38	76	11398	1.08	ug/L 91
28) methylene chloride	6.64	84	3966	1.04	ug/L 98
29) methyl acetate	6.46	74	448	0.95	ug/L # 1
31) methyl tert butyl ether	6.89	73	10313	1.14	ug/L 85
32) trans-1,2-dichloroethene	6.95	96	4367	1.17	ug/L 88
33) di-isopropyl ether	7.42	45	13716	0.98	ug/L 75
35) 1,1-dichloroethane	7.45	63	6746	1.11	ug/L 98
36) chloroprene	7.54	53	5523	1.17	ug/L 93
37) acrylonitrile	6.93	53	7280	6.58	ug/L 86
38) vinyl acetate	7.43	86	455	0.91	ug/L # 1
39) ethyl tert-butyl ether	7.83	59	11692	0.94	ug/L 94
41) 2,2-dichloropropane	8.11	77	3273	1.08	ug/L 82
42) cis-1,2-dichloroethene	8.11	96	4154	1.21	ug/L 83
44) propionitrile	8.19	54	5545	10.32	ug/L 97
45) bromochloromethane	8.40	128	1933	1.07	ug/L 91
46) tetrahydrofuran	8.43	42	1471	1.07	ug/L 97
47) chloroform	8.44	85	4552	1.13	ug/L 94
48) T-BUTYL FORMATE	8.46	59	2382	1.01	ug/L # 77
51) freon 113	5.97	151	2213	0.98	ug/L 91
52) methacrylonitrile	8.34	41	2492	1.03	ug/L 90
53) 1,1,1-trichloroethane	8.67	97	4307	1.07	ug/L 75

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54244.D
 Acq On : 5 Aug 2015 10:35 am
 Operator : TOANP
 Sample : ic2289-1
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Aug 10 08:18:11 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:57:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
54) cyclohexane	8.73	84	4874	1.12	ug/L #	74
57) epichlorohydrin	10.63	57	2058	5.33	ug/L	90
58) n-butyl alcohol	9.54	56	5784	44.30	ug/L	97
59) carbon tetrachloride	8.86	117	4461	1.11	ug/L	90
60) 1,1-dichloropropene	8.84	75	5058	1.20	ug/L	97
61) hexane	7.20	57	4109	1.19	ug/L	80
63) benzene	9.08	78	15599	1.09	ug/L	100
65) tert-amyl methyl ether	9.09	87	2145	0.91	ug/L #	81
66) heptane	9.21	57	2628	1.13	ug/L	94
68) 1,2-dichloroethane	9.10	62	5033	1.17	ug/L	94
69) trichloroethene	9.74	95	4058	1.19	ug/L	88
72) 2-nitropropane	10.50	41	1692	1.10	ug/L #	63
73) 2-chloroethyl vinyl ether	10.50	63	11570	5.35	ug/L	99
74) methyl methacrylate	9.99	100	864	1.01	ug/L #	44
75) 1,2-dichloropropane	10.01	63	4019	1.05	ug/L	93
76) dibromomethane	10.16	93	2526	1.11	ug/L	91
77) methylcyclohexane	9.94	83	4903	1.10	ug/L	96
78) bromodichloromethane	10.28	83	5244	1.13	ug/L	92
79) cis-1,3-dichloropropene	10.72	75	6576	1.10	ug/L	92
81) 4-methyl-2-pentanone	10.82	58	1463	1.03	ug/L	91
82) toluene	11.09	92	9587	1.18	ug/L	89
83) 3-methyl-1-butanol	10.84	55	3741	18.41	ug/L	83
84) trans-1,3-dichloropropene	11.30	75	5852	1.12	ug/L	93
85) ethyl methacrylate	11.27	69	4537	1.13	ug/L	84
86) 1,1,2-trichloroethane	11.52	83	3232	1.11	ug/L	98
87) 2-hexanone	11.70	58	1424	1.15	ug/L #	61
89) tetrachloroethene	11.70	164	4309	1.07	ug/L	97
90) 1,3-dichloropropane	11.71	76	6104	1.17	ug/L	97
91) butyl acetate	11.77	56	2584	0.92	ug/L #	84
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	4062	8.64	ug/L	92
93) dibromochloromethane	11.99	129	4516	1.11	ug/L	92
94) 1,2-dibromoethane	12.15	107	3864	1.16	ug/L	99
96) chlorobenzene	12.64	112	11061	1.17	ug/L	88
97) 1,1,1,2-tetrachloroethane	12.69	131	3990	1.17	ug/L	93
98) ethylbenzene	12.69	91	18635	1.18	ug/L	99
99) m,p-xylene	12.80	106	14263	2.41	ug/L	81
100) o-xylene	13.26	106	6866	1.15	ug/L	87
101) styrene	13.27	104	10951	1.14	ug/L	91
102) bromoform	13.57	173	3448	1.17	ug/L	93
104) isopropylbenzene	13.63	105	18302	1.22	ug/L	93
107) bromobenzene	14.07	156	5315	1.16	ug/L	94
108) 1,1,2,2-tetrachloroethane	13.98	83	4998	1.11	ug/L	87
109) trans-1,4-dichloro-2-butene	14.03	53	1326	1.10	ug/L	86
110) 1,2,3-trichloropropene	14.06	110	1091	0.97	ug/L	89
111) n-propylbenzene	14.08	91	20802	1.18	ug/L	96
113) 2-chlorotoluene	14.25	126	4472	1.19	ug/L #	77
114) 4-chlorotoluene	14.35	91	13383	1.15	ug/L	92
115) 1,3,5-trimethylbenzene	14.25	105	14996	1.18	ug/L	98
116) tert-butylbenzene	14.63	119	13117	1.15	ug/L	98
117) pentachloroethane	14.73	167	2665	1.07	ug/L	92

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54244.D
 Acq On : 5 Aug 2015 10:35 am
 Operator : TOANP
 Sample : ic2289-1
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Aug 10 08:18:11 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:57:32 2015
 Response via : Initial Calibration

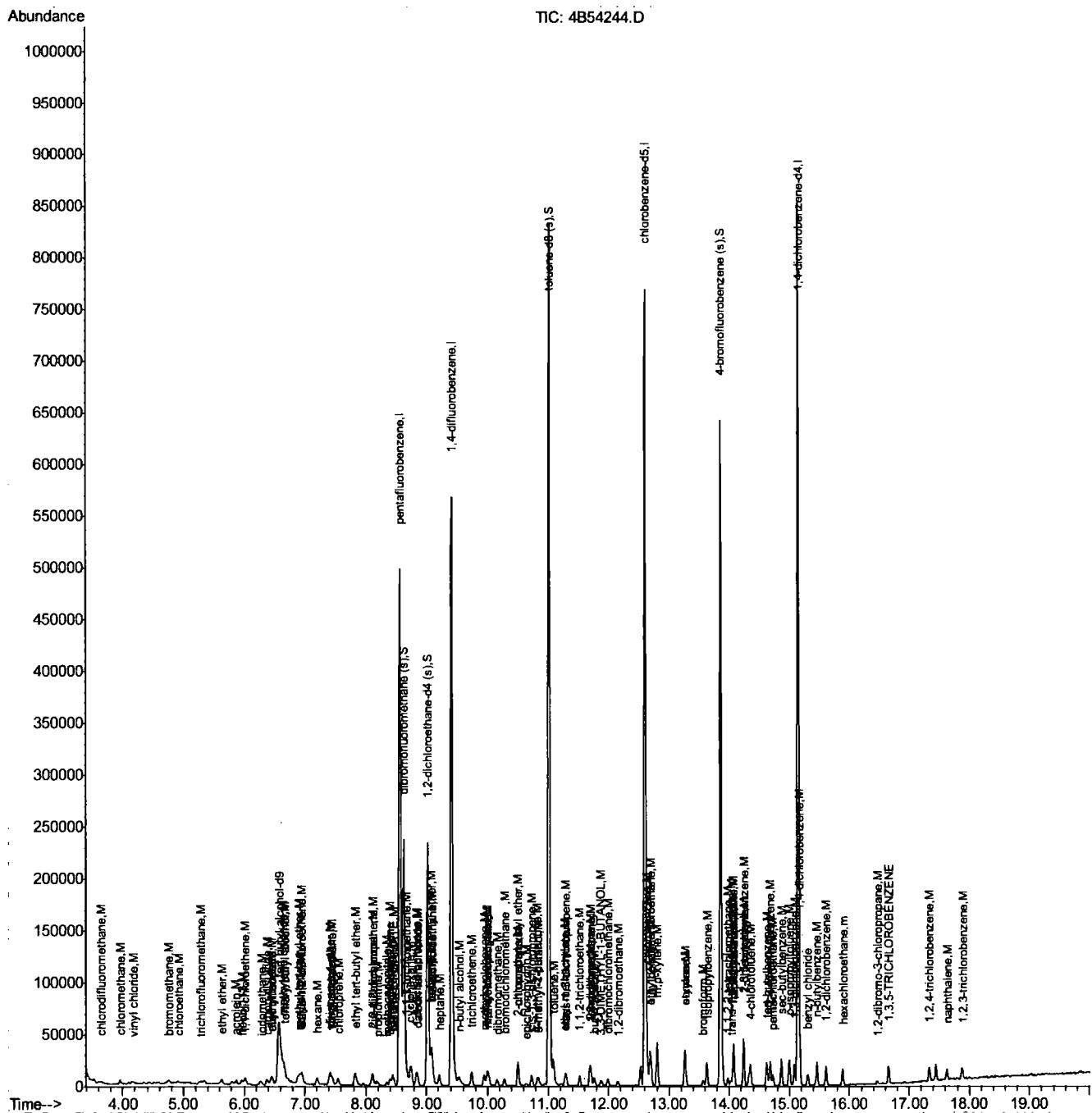
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
118) 1,2,4-trimethylbenzene	14.68	105	14265	1.12	ug/L	98
119) sec-butylbenzene	14.87	105	19440	1.18	ug/L	98
120) 1,3-dichlorobenzene	15.08	146	9710	1.12	ug/L	97
121) p-isopropyltoluene	15.00	119	15860	1.14	ug/L	97
122) 1,4-dichlorobenzene	15.17	146	9676	1.11	ug/L	92
123) benzyl chloride	15.30	91	9655	0.96	ug/L	#
124) 1,2-dichlorobenzene	15.61	146	9478	1.16	ug/L	91
126) n-butylbenzene	15.46	92	7635	1.16	ug/L	87
128) 1,2-dibromo-3-chloropropan	16.46	75	826	1.07	ug/L	#
129) 1,3,5-TRICHLOROBENZENE	16.65	180	6998	1.10	ug/L	99
130) 1,2,4-trichlorobenzene	17.33	180	5252	1.09	ug/L	96
132) naphthalene	17.62	128	8355	0.99	ug/L	98
133) 1,2,3-trichlorobenzene	17.88	180	4266	1.02	ug/L	93
134) hexachloroethane	15.89	201	3324	1.12	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
Data File : 4B54244.D
Acq On : 5 Aug 2015 10:35 am
Operator : TOANP
Sample : ic2289-1
Misc : MS88663,V4B2289,w,,,1
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Aug 10 08:18:11 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 13:57:32 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54245.D
 Acq On : 5 Aug 2015 11:03 am
 Operator : TOANP
 Sample : ic2289-2
 Misc : MS88663,V4B2289,w,,,
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 06 08:44:40 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:53:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	154019	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	458246	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	533965	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	468820	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	253142	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	159550	48.62	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	97.24%
50) 1,2-dichloroethane-d4 (s)	9.02	65	183532	50.40	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	100.80%
80) toluene-d8 (s)	11.02	98	629388	50.06	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	100.12%
105) 4-bromofluorobenzene (s)	13.86	95	222983	49.87	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.74%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.68	59	3724	9.82	ug/L # 69
3) 1,4-dioxane	10.12	88	1411	47.92	ug/L # 71
7) chlorodifluoromethane	3.65	51	9659	2.20	ug/L 95
8) dichlorodifluoromethane	3.62	85	8686	2.24	ug/L 85
10) chloromethane	3.96	52	3495	2.05	ug/L # 85
11) vinyl chloride	4.17	62	10299	2.25	ug/L 98
12) bromomethane	4.77	94	6188	1.92	ug/L 85
13) chloroethane	4.91	64	3927	2.03	ug/L 80
15) trichlorofluoromethane	5.28	101	9423	2.26	ug/L 93
19) ethyl ether	5.63	74	3631	2.11	ug/L # 73
21) acrolein	5.87	56	10833	19.45	ug/L 96
22) 1,1-dichloroethene	6.02	96	7880	2.13	ug/L 96
23) acetone	6.06	58	601	2.03	ug/L # 36
24) allyl chloride	6.46	76	4766	1.81	ug/L # 76
25) acetonitrile	6.45	40	5650	19.53	ug/L 86
26) iodomethane	6.28	142	14208	2.10	ug/L 99
27) carbon disulfide	6.38	76	22838	2.15	ug/L 97
28) methylene chloride	6.64	84	7908	2.07	ug/L 91
29) methyl acetate	6.46	74	940	1.99	ug/L # 33
30) 1-chloropropane	6.66	42	15074	2.01	ug/L 98
31) methyl tert butyl ether	6.90	73	19863	2.20	ug/L 90
32) trans-1,2-dichloroethene	6.96	96	7485	1.99	ug/L 96
33) di-isopropyl ether	7.41	45	27204	1.95	ug/L 97
35) 1,1-dichloroethane	7.45	63	14169	2.33	ug/L 93
36) chloroprene	7.55	53	10198	2.16	ug/L 95
37) acrylonitrile	6.93	53	12330	11.09	ug/L 91
38) vinyl acetate	7.44	86	1201	2.38	ug/L # 1
39) ethyl tert-butyl ether	7.83	59	24165	1.93	ug/L 96
40) ethyl acetate	8.09	45	1191	2.16	ug/L # 36
41) 2,2-dichloropropane	8.11	77	6636	2.18	ug/L 91
42) cis-1,2-dichloroethene	8.11	96	8200	2.39	ug/L 98
44) propionitrile	8.19	54	10716	19.86	ug/L 98
45) bromochloromethane	8.40	128	3878	2.15	ug/L 86

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54245.D
 Acq On : 5 Aug 2015 11:03 am
 Operator : TOANP
 Sample : ic2289-2
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 06 08:44:40 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:53:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) tetrahydrofuran	8.43	42	2774	2.02	ug/L	92
47) chloroform	8.44	85	8718	2.16	ug/L	95
48) T-BUTYL FORMATE	8.46	59	4749	2.01	ug/L	85
51) freon 113	5.96	151	4410	1.95	ug/L	88
52) methacrylonitrile	8.35	41	5052	2.08	ug/L	90
53) 1,1,1-trichloroethane	8.67	97	8807	2.18	ug/L	96
54) cyclohexane	8.74	84	9470	2.17	ug/L	97
57) epichlorohydrin	10.64	57	3465	8.87	ug/L	89
58) n-butyl alcohol	9.54	56	9630	72.95	ug/L	89
59) carbon tetrachloride	8.86	117	8815	2.17	ug/L	97
60) 1,1-dichloropropene	8.83	75	9953	2.34	ug/L	97
61) hexane	7.20	57	7648	2.18	ug/L	96
63) benzene	9.08	78	30380	2.11	ug/L	98
64) iso-octane	9.05	57	24406	2.36	ug/L	97
65) tert-amyl methyl ether	9.09	87	5012	2.11	ug/L	# 78
66) heptane	9.21	57	5509	2.35	ug/L	93
68) 1,2-dichloroethane	9.10	62	10057	2.31	ug/L	94
69) trichloroethene	9.74	95	7554	2.20	ug/L	97
71) ethyl acrylate	9.73	55	10852	2.18	ug/L	98
72) 2-nitropropane	10.51	41	3228	2.08	ug/L	# 80
73) 2-chloroethyl vinyl ether	10.50	63	23271	10.64	ug/L	96
74) methyl methacrylate	9.99	100	1658	1.92	ug/L	# 86
75) 1,2-dichloropropane	10.01	63	8626	2.22	ug/L	97
76) dibromomethane	10.16	93	4555	1.97	ug/L	84
77) methylcyclohexane	9.94	83	9847	2.19	ug/L	92
78) bromodichloromethane	10.27	83	10576	2.26	ug/L	92
79) cis-1,3-dichloropropene	10.72	75	13246	2.20	ug/L	95
81) 4-methyl-2-pentanone	10.82	58	2863	2.00	ug/L	# 78
82) toluene	11.09	92	18644	2.26	ug/L	97
83) 3-methyl-1-butanol	10.84	55	7245	35.27	ug/L	90
84) trans-1,3-dichloropropene	11.30	75	11430	2.16	ug/L	98
85) ethyl methacrylate	11.28	69	8731	2.15	ug/L	97
86) 1,1,2-trichloroethane	11.52	83	6418	2.17	ug/L	90
87) 2-hexanone	11.70	58	2491	1.99	ug/L	# 72
89) tetrachloroethene	11.69	164	8798	2.20	ug/L	86
90) 1,3-dichloropropane	11.71	76	11870	2.29	ug/L	97
91) butyl acetate	11.76	56	4842	1.73	ug/L	95
92) 3,3-DIMETHYL-1-BUTANOL	11.88	57	7873	16.79	ug/L	94
93) dibromochloromethane	11.99	129	8693	2.14	ug/L	98
94) 1,2-dibromoethane	12.15	107	7294	2.20	ug/L	93
96) chlorobenzene	12.64	112	21905	2.32	ug/L	91
97) 1,1,1,2-tetrachloroethane	12.70	131	7589	2.23	ug/L	99
98) ethylbenzene	12.68	91	35982	2.27	ug/L	98
99) m,p-xylene	12.80	106	26525	4.49	ug/L	95
100) o-xylene	13.25	106	13023	2.20	ug/L	94
101) styrene	13.27	104	20927	2.18	ug/L	97
102) bromoform	13.57	173	6228	2.12	ug/L	92
104) isopropylbenzene	13.63	105	33782	2.26	ug/L	96
106) cyclohexanone	13.83	55	9492	16.25	ug/L	98
107) bromobenzene	14.08	156	9969	2.19	ug/L	97

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54245.D
 Acq On : 5 Aug 2015 11:03 am
 Operator : TOANP
 Sample : ic2289-2
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 06 08:44:40 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:53:30 2015
 Response via : Initial Calibration

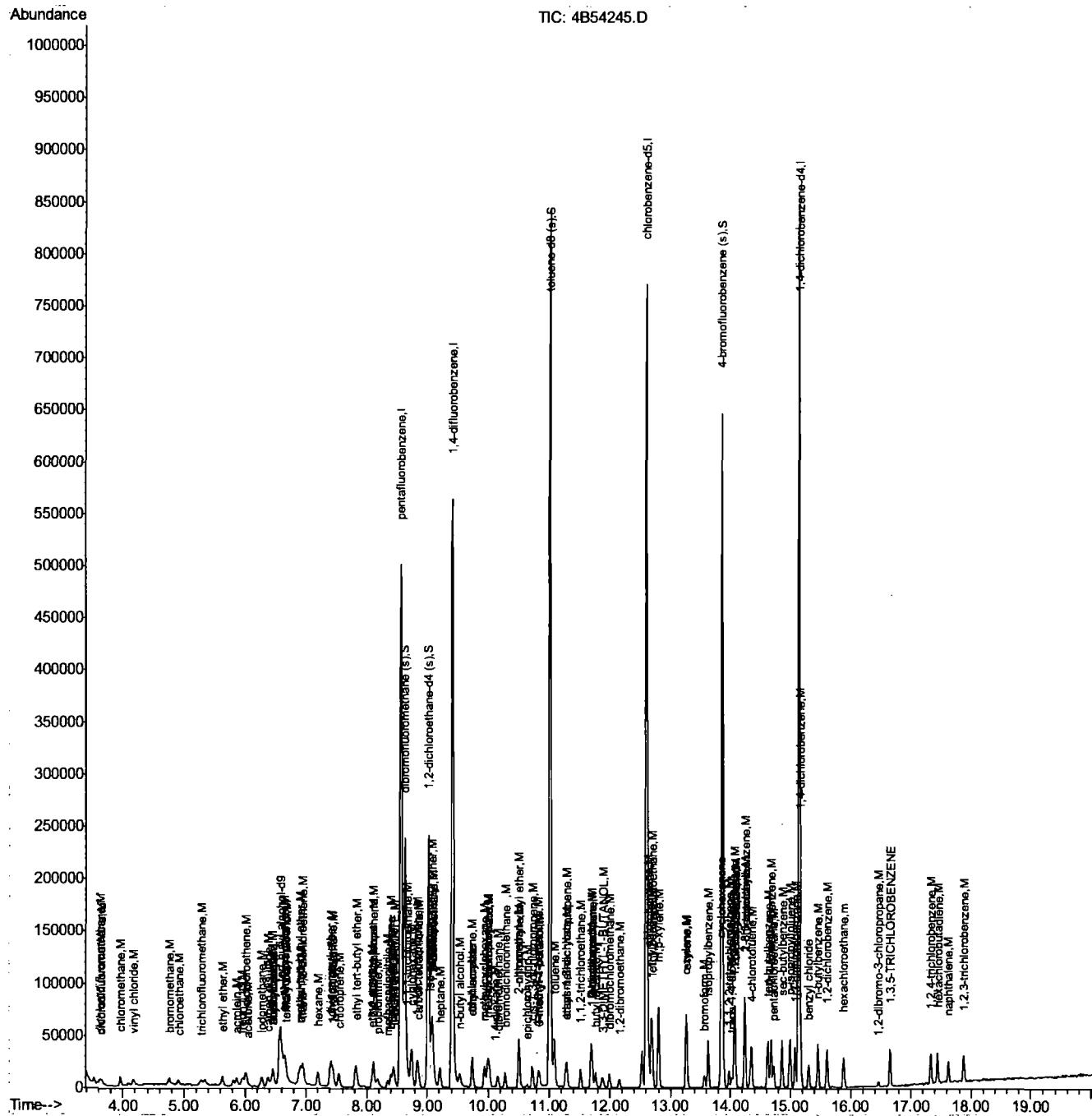
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
108) 1,1,2,2-tetrachloroethane	13.97	83	9934	2.21	ug/L	95
109) trans-1,4-dichloro-2-butene	14.02	53	2262	1.88	ug/L	91
110) 1,2,3-trichloropropane	14.06	110	2306	2.07	ug/L	92
111) n-propylbenzene	14.08	91	39209	2.23	ug/L	98
113) 2-chlorotoluene	14.24	126	8751	2.33	ug/L	91
114) 4-chlorotoluene	14.35	91	25519	2.21	ug/L	98
115) 1,3,5-trimethylbenzene	14.24	105	28842	2.28	ug/L	98
116) tert-butylbenzene	14.62	119	24982	2.19	ug/L	97
117) pentachloroethane	14.72	167	5440	2.20	ug/L	92
118) 1,2,4-trimethylbenzene	14.68	105	28044	2.22	ug/L	98
119) sec-butylbenzene	14.87	105	37382	2.27	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	18557	2.16	ug/L	93
121) p-isopropyltoluene	15.00	119	30933	2.24	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	18638	2.15	ug/L	96
123) benzyl chloride	15.30	91	17614	1.76	ug/L	99
124) 1,2-dichlorobenzene	15.61	146	17338	2.13	ug/L	97
126) n-butylbenzene	15.46	92	14589	2.23	ug/L	95
128) 1,2-dibromo-3-chloropropan	16.47	75	1478	1.92	ug/L	79
129) 1,3,5-TRICHLOROBENZENE	16.65	180	14190	2.25	ug/L	95
130) 1,2,4-trichlorobenzene	17.33	180	10099	2.11	ug/L	90
131) hexachlorobutadiene	17.44	225	7220	2.16	ug/L	93
132) naphthalene	17.62	128	16981	2.03	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	8562	2.06	ug/L	90
134) hexachloroethane	15.89	201	6297	2.13	ug/L	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54245.D
Acq On : 5 Aug 2015 11:03 am
Operator : TOANP
Sample : ic2289-2
Misc : MS88663,V4B2289,w,,,,1
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Aug 06 08:44:40 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 13:53:30 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54246.D
 Acq On : 5 Aug 2015 11:31 am
 Operator : TOANP
 Sample : ic2289-5
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 05 13:55:02 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:53:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	162164	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	454275	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	525775	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	462724	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	255087	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	163731	50.33	ug/L	0.00
Spiked Amount 50.000	Range 76 - 120		Recovery	=	100.66%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	182681	50.60	ug/L	0.00
Spiked Amount 50.000	Range 73 - 122		Recovery	=	101.20%	
80) toluene-d8 (s)	11.02	98	616561	49.80	ug/L	0.00
Spiked Amount 50.000	Range 84 - 119		Recovery	=	99.60%	
105) 4-bromofluorobenzene (s)	13.86	95	224431	49.81	ug/L	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery	=	99.62%	

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.67	59	9342	23.39	ug/L
3) 1,4-dioxane	10.12	88	4284	138.18	ug/L
7) chlorodifluoromethane	3.64	51	21460	4.93	ug/L
8) dichlorodifluoromethane	3.61	85	22589	5.88	ug/L
10) chloromethane	3.96	52	8543	5.05	ug/L
11) vinyl chloride	4.17	62	24768	5.46	ug/L
12) bromomethane	4.75	94	14298	4.48	ug/L
13) chloroethane	4.91	64	9996	5.22	ug/L
15) trichlorofluoromethane	5.28	101	23817	5.76	ug/L
19) ethyl ether	5.62	74	8326	4.87	ug/L
20) 2-chloropropane	5.81	39	3012	4.32	ug/L
21) acrolein	5.87	56	27925	50.57	ug/L
22) 1,1-dichloroethene	6.02	96	17826	4.86	ug/L
23) acetone	6.05	58	1487	5.08	ug/L #
24) allyl chloride	6.46	76	11576	4.42	ug/L
25) acetonitrile	6.45	40	13312	46.42	ug/L
26) iodomethane	6.28	142	33601	5.01	ug/L
27) carbon disulfide	6.38	76	51918	4.92	ug/L
28) methylene chloride	6.64	84	18751	4.95	ug/L
29) methyl acetate	6.44	74	2176	4.65	ug/L #
30) 1-chloropropane	6.66	42	34118	4.59	ug/L
31) methyl tert butyl ether	6.89	73	48034	5.36	ug/L
32) trans-1,2-dichloroethene	6.95	96	17388	4.67	ug/L
33) di-isopropyl ether	7.41	45	61480	4.44	ug/L
34) 2-butanone	8.08	72	1802	5.51	ug/L #
35) 1,1-dichloroethane	7.45	63	33333	5.53	ug/L
36) chloroprene	7.55	53	24327	5.19	ug/L
37) acrylonitrile	6.93	53	32842	29.81	ug/L
38) vinyl acetate	7.43	86	2329	4.66	ug/L
39) ethyl tert-butyl ether	7.83	59	53574	4.32	ug/L
40) ethyl acetate	8.09	45	2588	4.74	ug/L
41) 2,2-dichloropropane	8.11	77	14526	4.81	ug/L
42) cis-1,2-dichloroethene	8.11	96	18214	5.34	ug/L

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54246.D
 Acq On : 5 Aug 2015 11:31 am
 Operator : TOANP
 Sample : ic2289-5
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 05 13:55:02 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:53:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.17	85	2310	5.02	ug/L #	65
44) propionitrile	8.19	54	27841	52.04	ug/L	97
45) bromochloromethane	8.40	128	9368	5.23	ug/L	94
46) tetrahydrofuran	8.43	42	6501	4.77	ug/L	92
47) chloroform	8.44	85	19751	4.93	ug/L	97
48) T-BUTYL FORMATE	8.47	59	11385	4.87	ug/L	88
51) freon 113	5.96	151	11367	5.06	ug/L	96
52) methacrylonitrile	8.35	41	11856	4.92	ug/L	93
53) 1,1,1-trichloroethane	8.67	97	20376	5.08	ug/L	98
54) cyclohexane	8.74	84	22770	5.26	ug/L	95
55) iso-butyl alcohol	8.83	43	8121	47.70	ug/L	94
57) epichlorohydrin	10.64	57	9359	24.34	ug/L	98
58) n-butyl alcohol	9.53	56	26854	206.60	ug/L	95
59) carbon tetrachloride	8.86	117	20465	5.12	ug/L	94
60) 1,1-dichloropropene	8.83	75	23325	5.57	ug/L	96
61) hexane	7.20	57	18603	5.40	ug/L	94
63) benzene	9.08	78	68859	4.85	ug/L	98
64) iso-octane	9.07	57	59474	5.84	ug/L	98
65) tert-amyl methyl ether	9.09	87	10708	4.58	ug/L #	86
66) heptane	9.21	57	12942	5.60	ug/L	96
67) isopropyl acetate	8.99	61	10596	2.65	ug/L #	67
68) 1,2-dichloroethane	9.10	62	23204	5.41	ug/L	99
69) trichloroethene	9.74	95	18278	5.40	ug/L	99
71) ethyl acrylate	9.73	55	32881	6.71	ug/L	96
72) 2-nitropropane	10.50	41	7628	5.00	ug/L	96
73) 2-chloroethyl vinyl ether	10.50	63	54357	25.23	ug/L	98
74) methyl methacrylate	9.99	100	4362	5.12	ug/L #	66
75) 1,2-dichloropropane	10.00	63	19175	5.01	ug/L	99
76) dibromomethane	10.16	93	11447	5.04	ug/L	91
77) methylcyclohexane	9.95	83	24696	5.58	ug/L	94
78) bromodichloromethane	10.28	83	25253	5.48	ug/L	95
79) cis-1,3-dichloropropene	10.73	75	31392	5.30	ug/L	98
81) 4-methyl-2-pentanone	10.82	58	7935	5.62	ug/L #	81
82) toluene	11.09	92	43511	5.37	ug/L	99
83) 3-methyl-1-butanol	10.83	55	17788	87.95	ug/L	98
84) trans-1,3-dichloropropene	11.30	75	28985	5.56	ug/L	90
85) ethyl methacrylate	11.28	69	22111	5.54	ug/L	94
86) 1,1,2-trichloroethane	11.52	83	14633	5.03	ug/L	99
87) 2-hexanone	11.70	58	7235	5.86	ug/L	84
89) tetrachloroethene	11.69	164	20781	5.26	ug/L	93
90) 1,3-dichloropropane	11.71	76	28621	5.59	ug/L	96
91) butyl acetate	11.76	56	12776	4.63	ug/L	94
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	19357	41.84	ug/L	97
93) dibromochloromethane	11.99	129	20497	5.12	ug/L	99
94) 1,2-dibromoethane	12.15	107	17981	5.49	ug/L	99
96) chlorobenzene	12.64	112	48909	5.24	ug/L	97
97) 1,1,1,2-tetrachloroethane	12.70	131	18650	5.55	ug/L	94
98) ethylbenzene	12.69	91	83346	5.34	ug/L	96
99) m,p-xylene	12.80	106	62446	10.70	ug/L	100
100) o-xylene	13.25	106	31758	5.43	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54246.D
 Acq On : 5 Aug 2015 11:31 am
 Operator : TOANP
 Sample : ic2289-5
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 05 13:55:02 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:53:30 2015
 Response via : Initial Calibration

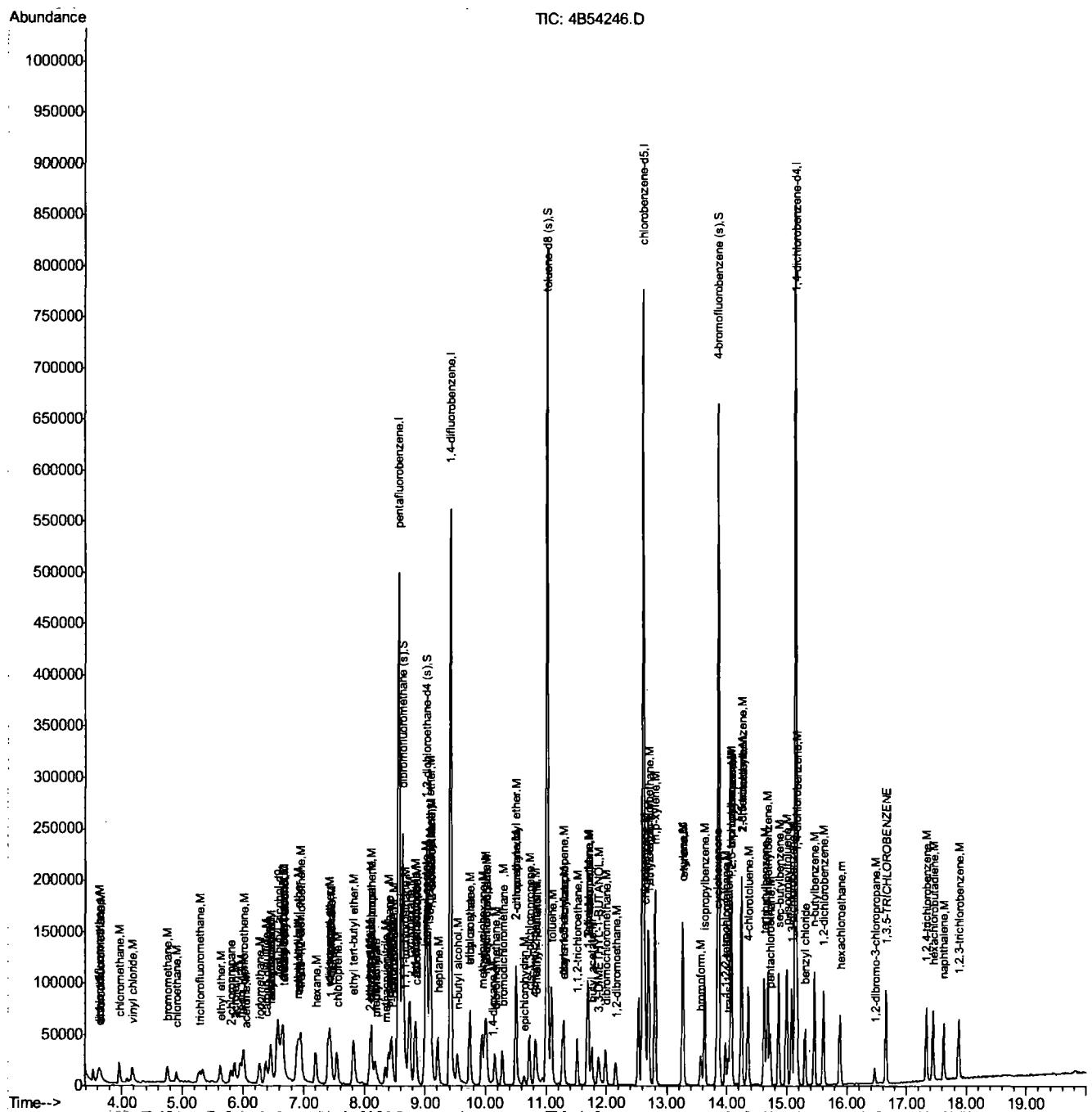
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
101) styrene	13.27	104	51009	5.40	ug/L	96
102) bromoform	13.57	173	15914	5.49	ug/L	98
104) isopropylbenzene	13.63	105	81026	5.38	ug/L	99
106) cyclohexanone	13.83	55	26972	45.84	ug/L	96
107) bromobenzene	14.08	156	24320	5.29	ug/L	99
108) 1,1,2,2-tetrachloroethane	13.98	83	24652	5.45	ug/L	95
109) trans-1,4-dichloro-2-buten	14.02	53	6398	5.29	ug/L	97
110) 1,2,3-trichloropropane	14.06	110	6089	5.43	ug/L	96
111) n-propylbenzene	14.08	91	94454	5.32	ug/L	98
113) 2-chlorotoluene	14.24	126	20347	5.38	ug/L	100
114) 4-chlorotoluene	14.35	91	60152	5.16	ug/L	97
115) 1,3,5-trimethylbenzene	14.24	105	67454	5.28	ug/L	98
116) tert-butylbenzene	14.63	119	59531	5.19	ug/L	99
117) pentachloroethane	14.72	167	12451	4.99	ug/L	98
118) 1,2,4-trimethylbenzene	14.68	105	67240	5.29	ug/L	98
119) sec-butylbenzene	14.87	105	89332	5.39	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	44161	5.10	ug/L	97
121) p-isopropyltoluene	15.00	119	75077	5.40	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	43088	4.94	ug/L	98
123) benzyl chloride	15.30	91	43407	4.30	ug/L	98
124) 1,2-dichlorobenzene	15.61	146	42296	5.15	ug/L	98
126) n-butylbenzene	15.46	92	36109	5.47	ug/L	97
128) 1,2-dibromo-3-chloropropan	16.46	75	3677	4.75	ug/L	82
129) 1,3,5-TRICHLOROBENZENE	16.65	180	33747	5.30	ug/L	98
130) 1,2,4-trichlorobenzene	17.33	180	25250	5.22	ug/L	94
131) hexachlorobutadiene	17.44	225	17204	5.11	ug/L	97
132) naphthalene	17.62	128	43667	5.17	ug/L	100
133) 1,2,3-trichlorobenzene	17.88	180	21445	5.12	ug/L	97
134) hexachloroethane	15.89	201	15447	5.20	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
Data File : 4B54246.D
Acq On : 5 Aug 2015 11:31 am
Operator : TOANP
Sample : ic2289-5
Misc : MS88663,V4B2289,w,,,1
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Aug 05 13:55:02 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 13:53:30 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54247.D
 Acq On : 5 Aug 2015 11:59 am
 Operator : TOANP
 Sample : ic2289-10
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Aug 05 13:46:01 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:44:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	156272	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	441872	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	519156	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	463675	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	259714	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	161777	51.12	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	102.24%
50) 1,2-dichloroethane-d4 (s)	9.02	65	176378	50.23	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	100.46%
80) toluene-d8 (s)	11.02	98	615564	50.36	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	100.72%
105) 4-bromofluorobenzene (s)	13.86	95	225049	49.06	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.12%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.68	59	19388	50.38	ug/L
3) 1,4-dioxane	10.12	88	8688	290.80	ug/L
7) chlorodifluoromethane	3.64	51	43835	10.36	ug/L
8) dichlorodifluoromethane	3.61	85	44521	11.91	ug/L
10) chloromethane	3.96	52	16125	9.80	ug/L
11) vinyl chloride	4.17	62	48692	11.04	ug/L
12) bromomethane	4.76	94	27980	9.02	ug/L
13) chloroethane	4.91	64	19384	10.41	ug/L
15) trichlorofluoromethane	5.28	101	46904	11.66	ug/L
19) ethyl ether	5.63	74	17070	10.27	ug/L
20) 2-chloropropane	5.82	39	6493	9.58	ug/L
21) acrolein	5.86	56	57328	106.72	ug/L
22) 1,1-dichloroethene	6.01	96	36664	10.28	ug/L
23) acetone	6.06	58	2758	9.68	ug/L #
24) allyl chloride	6.46	76	28061	11.02	ug/L #
25) acetonitrile	6.45	40	30542	109.50	ug/L
26) iodomethane	6.27	142	70690	10.83	ug/L
27) carbon disulfide	6.38	76	109861	10.71	ug/L
28) methylene chloride	6.65	84	39476	10.72	ug/L
29) methyl acetate	6.44	74	5110	11.22	ug/L #
30) 1-chloropropane	6.66	42	66479	9.20	ug/L
31) methyl tert butyl ether	6.89	73	96930	11.11	ug/L
32) trans-1,2-dichloroethene	6.95	96	35968	9.94	ug/L
33) di-isopropyl ether	7.41	45	125477	9.31	ug/L
34) 2-butanone	8.09	72	3814	12.00	ug/L #
35) 1,1-dichloroethane	7.45	63	68554	11.70	ug/L
36) chloroprene	7.55	53	48716	10.69	ug/L
37) acrylonitrile	6.92	53	66799	62.33	ug/L
38) vinyl acetate	7.43	86	4739	9.74	ug/L
39) ethyl tert-butyl ether	7.83	59	109885	9.11	ug/L
40) ethyl acetate	8.09	45	5446	10.25	ug/L
41) 2,2-dichloropropane	8.11	77	29884	10.17	ug/L
42) cis-1,2-dichloroethene	8.11	96	38288	11.55	ug/L

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54247.D
 Acq On : 5 Aug 2015 11:59 am
 Operator : TOANP
 Sample : ic2289-10
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Aug 05 13:46:01 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:44:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.17	85	4945	11.04	ug/L	# 86
44) propionitrile	8.19	54	55994	107.59	ug/L	91
45) bromochloromethane	8.40	128	19683	11.29	ug/L	92
46) tetrahydrofuran	8.43	42	12786	9.64	ug/L	97
47) chloroform	8.44	85	41508	10.66	ug/L	96
48) T-BUTYL FORMATE	8.47	59	23667	10.40	ug/L	96
51) freon 113	5.96	151	22630	10.35	ug/L	96
52) methacrylonitrile	8.34	41	24067	10.27	ug/L	94
53) 1,1,1-trichloroethane	8.67	97	43974	11.27	ug/L	96
54) cyclohexane	8.74	84	47618	11.31	ug/L	# 82
55) iso-butyl alcohol	8.84	43	14607	88.20	ug/L	99
57) epichlorohydrin	10.64	57	18910	49.80	ug/L	95
58) n-butyl alcohol	9.53	56	51734	403.09	ug/L	99
59) carbon tetrachloride	8.86	117	44023	11.15	ug/L	95
60) 1,1-dichloropropene	8.84	75	49092	11.87	ug/L	98
61) hexane	7.20	57	36571	10.75	ug/L	97
63) benzene	9.08	78	143094	10.21	ug/L	100
64) iso-octane	9.07	57	114940	11.43	ug/L	98
65) tert-amyl methyl ether	9.09	87	22078	9.57	ug/L	94
66) heptane	9.21	57	23430	10.27	ug/L	96
67) isopropyl acetate	8.99	61	17962	4.55	ug/L	92
68) 1,2-dichloroethane	9.10	62	48963	11.56	ug/L	97
69) trichloroethene	9.74	95	38485	11.51	ug/L	96
71) ethyl acrylate	9.73	55	52003	10.75	ug/L	100
72) 2-nitropropane	10.50	41	15536	10.31	ug/L	# 85
73) 2-chloroethyl vinyl ether	10.50	63	113330	53.28	ug/L	99
74) methyl methacrylate	9.99	100	8470	10.07	ug/L	# 92
75) 1,2-dichloropropane	10.00	63	41602	11.01	ug/L	97
76) dibromomethane	10.16	93	24416	10.89	ug/L	97
77) methylcyclohexane	9.94	83	48507	11.10	ug/L	96
78) bromodichloromethane	10.28	83	51739	11.37	ug/L	96
79) cis-1,3-dichloropropene	10.73	75	65515	11.19	ug/L	98
81) 4-methyl-2-pentanone	10.82	58	15353	11.01	ug/L	# 84
82) toluene	11.09	92	89345	11.16	ug/L	99
83) 3-methyl-1-butanol	10.83	55	33955	170.02	ug/L	98
84) trans-1,3-dichloropropene	11.29	75	58598	11.39	ug/L	97
85) ethyl methacrylate	11.28	69	46162	11.71	ug/L	99
86) 1,1,2-trichloroethane	11.52	83	30525	10.63	ug/L	98
87) 2-hexanone	11.69	58	13613	11.16	ug/L	96
89) tetrachloroethene	11.70	164	42406	10.72	ug/L	92
90) 1,3-dichloropropane	11.71	76	57169	11.14	ug/L	97
91) butyl acetate	11.76	56	25650	9.28	ug/L	89
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	37965	81.88	ug/L	99
93) dibromochloromethane	11.99	129	43061	10.73	ug/L	97
94) 1,2-dibromoethane	12.15	107	36547	11.14	ug/L	99
96) chlorobenzene	12.63	112	102463	10.96	ug/L	99
97) 1,1,1,2-tetrachloroethane	12.70	131	39056	11.60	ug/L	96
98) ethylbenzene	12.69	91	172101	11.00	ug/L	100
99) m,p-xylene	12.80	106	130958	22.40	ug/L	99
100) o-xylene	13.26	106	65798	11.22	ug/L	100

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54247.D
 Acq On : 5 Aug 2015 11:59 am
 Operator : TOANP
 Sample : ic2289-10
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Aug 05 13:46:01 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:44:20 2015
 Response via : Initial Calibration

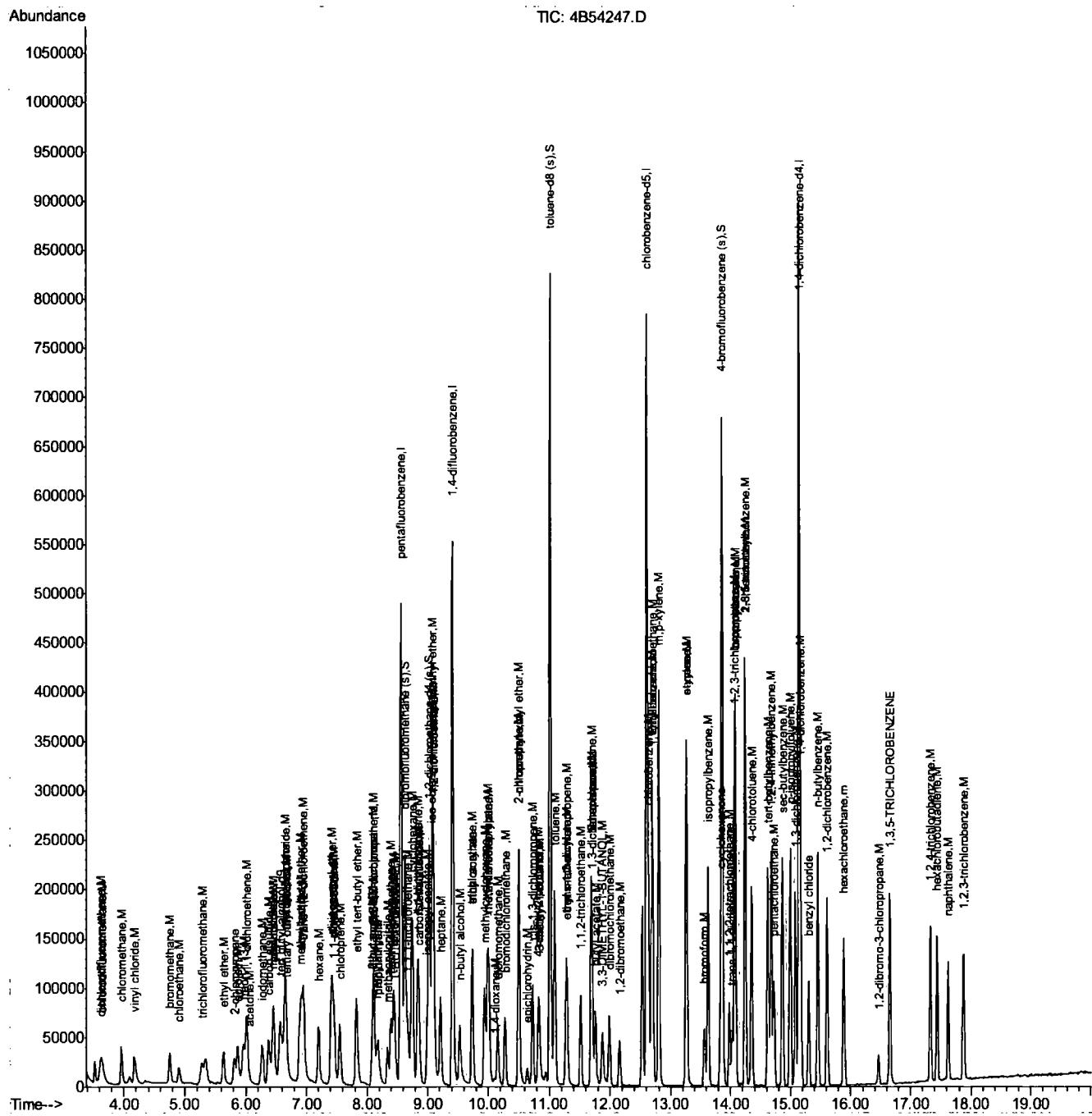
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
101) styrene	13.27	104	109974	11.61	ug/L	96
102) bromoform	13.57	173	32898	11.33	ug/L	94
104) isopropylbenzene	13.63	105	172120	11.22	ug/L	99
106) cyclohexanone	13.83	55	51662	86.23	ug/L	99
107) bromobenzene	14.07	156	50522	10.80	ug/L	91
108) 1,1,2,2-tetrachloroethane	13.98	83	50647	11.00	ug/L	98
109) trans-1,4-dichloro-2-buten	14.02	53	13345	10.83	ug/L	93
110) 1,2,3-trichloropropane	14.06	110	12435	10.89	ug/L	99
111) n-propylbenzene	14.08	91	198532	10.99	ug/L	99
113) 2-chlorotoluene	14.24	126	44347	11.52	ug/L	97
114) 4-chlorotoluene	14.35	91	127261	10.73	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	144636	11.12	ug/L	100
116) tert-butylbenzene	14.62	119	126427	10.82	ug/L	99
117) pentachloroethane	14.72	167	25449	10.02	ug/L	96
118) 1,2,4-trimethylbenzene	14.68	105	145042	11.20	ug/L	98
119) sec-butylbenzene	14.87	105	191619	11.36	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	95046	10.78	ug/L	99
121) p-isopropyltoluene	15.00	119	159749	11.29	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	92282	10.39	ug/L	100
123) benzyl chloride	15.30	91	85386	8.30	ug/L	99
124) 1,2-dichlorobenzene	15.61	146	91280	10.92	ug/L	98
126) n-butylbenzene	15.46	92	80220	11.95	ug/L	96
128) 1,2-dibromo-3-chloropropan	16.47	75	8051	10.21	ug/L	80
129) 1,3,5-TRICHLOROBENZENE	16.65	180	72920	11.25	ug/L	99
130) 1,2,4-trichlorobenzene	17.33	180	55206	11.22	ug/L	97
131) hexachlorobutadiene	17.44	225	37662	10.99	ug/L	99
132) naphthalene	17.62	128	95648	11.13	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	46835	10.99	ug/L	98
134) hexachloroethane	15.89	201	33253	10.98	ug/L	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54247.D
 Acq On : 5 Aug 2015 11:59 am
 Operator : TOANP
 Sample : ic2289-10
 Misc : MS88663, V4B2289, w, , , 1
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Aug 05 13:46:01 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:44:20 2015
 Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54248.D
 Acq On : 5 Aug 2015 12:27 pm
 Operator : TOANP
 Sample : ic2289-20
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Aug 05 13:44:07 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:42:40 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.57	65	149949	500.00	ug/L	-0.01
5) pentafluorobenzene	8.56	168	438540	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	518317	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	458544	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	262344	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	161373	51.38	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	102.76%
50) 1,2-dichloroethane-d4 (s)	9.02	65	177356	50.89	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	101.78%
80) toluene-d8 (s)	11.02	98	609847	49.97	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	99.94%
105) 4-bromofluorobenzene (s)	13.86	95	227455	49.08	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.16%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.67	59	37464	101.45	ug/L
3) 1,4-dioxane	10.11	88	17935	625.63	ug/L
7) chlorodifluoromethane	3.64	51	85293	20.30	ug/L
8) dichlorodifluoromethane	3.61	85	80693	21.75	ug/L
10) chloromethane	3.96	52	30540	18.70	ug/L
11) vinyl chloride	4.17	62	90837	20.75	ug/L
12) bromomethane	4.76	94	50648	16.46	ug/L
13) chloroethane	4.91	64	36225	19.61	ug/L
15) trichlorofluoromethane	5.28	101	86196	21.60	ug/L
19) ethyl ether	5.63	74	32937	19.96	ug/L
20) 2-chloropropane	5.80	39	11535	17.15	ug/L
21) acrolein	5.87	56	109348	205.11	ug/L
22) 1,1-dichloroethene	6.02	96	66579	18.81	ug/L
23) acetone	6.06	58	5809	20.54	ug/L
24) allyl chloride	6.46	76	45175	17.88	ug/L
25) acetonitrile	6.46	40	58922	212.85	ug/L #
26) iodomethane	6.28	142	127887	19.74	ug/L
27) carbon disulfide	6.38	76	203132	19.95	ug/L
28) methylene chloride	6.65	84	71427	19.54	ug/L
29) methyl acetate	6.44	74	10394	23.00	ug/L
30) 1-chloropropane	6.66	42	117234	16.35	ug/L
31) methyl tert butyl ether	6.89	73	182935	21.14	ug/L
32) trans-1,2-dichloroethene	6.95	96	62927	17.52	ug/L
33) di-isopropyl ether	7.41	45	240795	18.00	ug/L
34) 2-butanone	8.08	72	7779	24.65	ug/L #
35) 1,1-dichloroethane	7.45	63	125912	21.65	ug/L
36) chloroprene	7.55	53	92024	20.34	ug/L
37) acrylonitrile	6.92	53	128033	120.37	ug/L
38) vinyl acetate	7.43	86	10022	20.75	ug/L
39) ethyl tert-butyl ether	7.83	59	213311	17.82	ug/L
40) ethyl acetate	8.09	45	10168	19.28	ug/L
41) 2,2-dichloropropane	8.11	77	52937	18.15	ug/L
42) cis-1,2-dichloroethene	8.11	96	69985	21.27	ug/L

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54248.D
 Acq On : 5 Aug 2015 12:27 pm
 Operator : TOANP
 Sample : ic2289-20
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Aug 05 13:44:07 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:42:40 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.16	85	9823	22.09	ug/L	95
44) propionitrile	8.19	54	105304	203.88	ug/L	96
45) bromochloromethane	8.40	128	37367	21.60	ug/L	98
46) tetrahydrofuran	8.43	42	24872	18.90	ug/L	94
47) chloroform	8.44	85	74533	19.28	ug/L	98
48) T-BUTYL FORMATE	8.46	59	49010	21.70	ug/L	97
51) freon 113	5.97	151	41465	19.11	ug/L	94
52) methacrylonitrile	8.34	41	45468	19.56	ug/L	94
53) 1,1,1-trichloroethane	8.67	97	79828	20.62	ug/L	98
54) cyclohexane	8.74	84	84445	20.21	ug/L	98
55) iso-butyl alcohol	8.83	43	27489	167.25	ug/L	99
57) epichlorohydrin	10.64	57	32595	85.98	ug/L	96
58) n-butyl alcohol	9.53	56	98594	769.45	ug/L	97
59) carbon tetrachloride	8.86	117	80738	20.48	ug/L	99
60) 1,1-dichloropropene	8.83	75	89819	21.75	ug/L	98
61) hexane	7.20	57	65776	19.36	ug/L	99
63) benzene	9.08	78	259291	18.54	ug/L	99
64) iso-octane	9.06	57	210051	20.93	ug/L	100
65) tert-amyl methyl ether	9.09	87	42400	18.41	ug/L	99
66) heptane	9.21	57	46675	20.48	ug/L	98
67) isopropyl acetate	8.99	61	30472	7.73	ug/L	97
68) 1,2-dichloroethane	9.10	62	90113	21.31	ug/L	99
69) trichloroethene	9.74	95	69249	20.75	ug/L	98
71) ethyl acrylate	9.73	55	94791	19.62	ug/L	100
72) 2-nitropropane	10.50	41	28999	19.28	ug/L	# 88
73) 2-chloroethyl vinyl ether	10.50	63	218618	102.95	ug/L	99
74) methyl methacrylate	9.99	100	17369	20.69	ug/L	96
75) 1,2-dichloropropane	10.00	63	76735	20.34	ug/L	100
76) dibromomethane	10.16	93	44981	20.09	ug/L	98
77) methylcyclohexane	9.94	83	90932	20.85	ug/L	99
78) bromodichloromethane	10.28	83	96423	21.23	ug/L	98
79) cis-1,3-dichloropropene	10.72	75	122438	20.95	ug/L	99
81) 4-methyl-2-pentanone	10.82	58	30533	21.92	ug/L	97
82) toluene	11.09	92	162561	20.34	ug/L	97
83) 3-methyl-1-butanol	10.83	55	65120	326.60	ug/L	98
84) trans-1,3-dichloropropene	11.29	75	111342	21.68	ug/L	97
85) ethyl methacrylate	11.27	69	88641	22.53	ug/L	99
86) 1,1,2-trichloroethane	11.52	83	56822	19.82	ug/L	98
87) 2-hexanone	11.69	58	26570	21.82	ug/L	91
89) tetrachloroethene	11.69	164	76677	19.59	ug/L	98
90) 1,3-dichloropropane	11.71	76	106519	20.99	ug/L	98
91) butyl acetate	11.76	56	47568	17.40	ug/L	89
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	72862	158.91	ug/L	99
93) dibromochloromethane	11.99	129	80482	20.28	ug/L	97
94) 1,2-dibromoethane	12.15	107	69496	21.41	ug/L	100
96) chlorobenzene	12.64	112	188554	20.39	ug/L	99
97) 1,1,1,2-tetrachloroethane	12.70	131	71720	21.54	ug/L	98
98) ethylbenzene	12.69	91	315714	20.41	ug/L	99
99) m,p-xylene	12.80	106	239973	41.51	ug/L	98
100) o-xylene	13.26	106	122336	21.09	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54248.D
 Acq On : 5 Aug 2015 12:27 pm
 Operator : TOANP
 Sample : ic2289-20
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 8 Sample Multiplier: 1

Quant Time: Aug 05 13:44:07 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 13:42:40 2015
 Response via : Initial Calibration

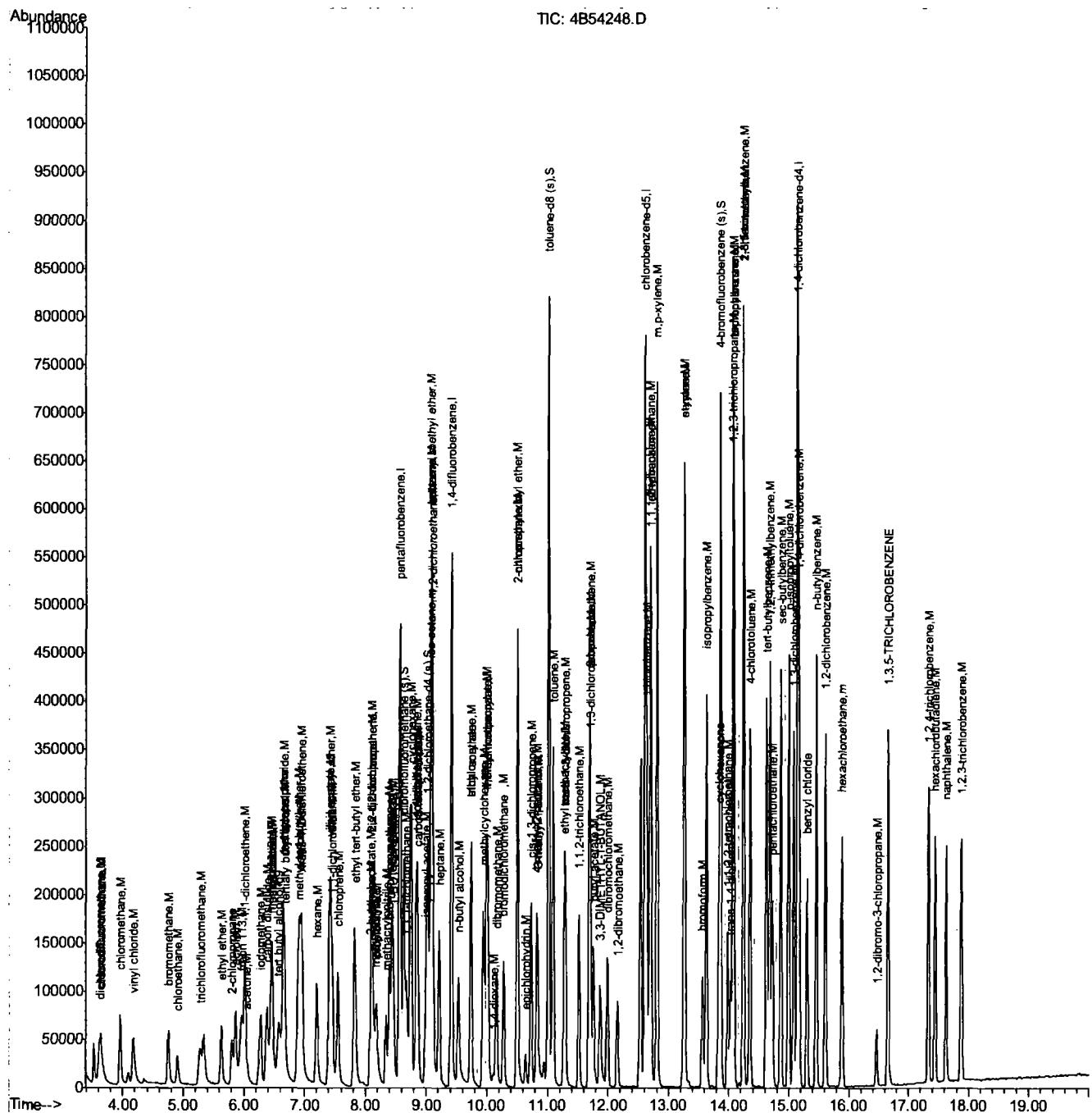
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
101) styrene	13.27	104	206596	22.05	ug/L	99
102) bromoform	13.57	173	63173	22.00	ug/L	97
104) isopropylbenzene	13.63	105	312620	20.18	ug/L	99
106) cyclohexanone	13.83	55	96749	159.87	ug/L	98
107) bromobenzene	14.07	156	94149	19.92	ug/L	95
108) 1,1,2,2-tetrachloroethane	13.98	83	95080	20.45	ug/L	99
109) trans-1,4-dichloro-2-buten	14.02	53	24658	19.82	ug/L	95
110) 1,2,3-trichloropropane	14.06	110	23951	20.76	ug/L	98
111) n-propylbenzene	14.08	91	359488	19.70	ug/L	99
113) 2-chlorotoluene	14.24	126	80714	20.77	ug/L	98
114) 4-chlorotoluene	14.35	91	232703	19.42	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	264712	20.15	ug/L	99
116) tert-butylbenzene	14.62	119	231944	19.64	ug/L	99
117) pentachloroethane	14.72	167	47764	18.62	ug/L	98
118) 1,2,4-trimethylbenzene	14.68	105	263513	20.14	ug/L	99
119) sec-butylbenzene	14.87	105	350667	20.58	ug/L	98
120) 1,3-dichlorobenzene	15.08	146	174107	19.55	ug/L	99
121) p-isopropyltoluene	15.00	119	296957	20.77	ug/L	98
122) 1,4-dichlorobenzene	15.17	146	171579	19.13	ug/L	97
123) benzyl chloride	15.30	91	168222	16.20	ug/L	100
124) 1,2-dichlorobenzene	15.61	146	171237	20.29	ug/L	99
126) n-butylbenzene	15.46	92	149839	22.09	ug/L	99
128) 1,2-dibromo-3-chloropropan	16.46	75	15204	19.08	ug/L	84
129) 1,3,5-TRICHLOROBENZENE	16.65	180	138324	21.13	ug/L	100
130) 1,2,4-trichlorobenzene	17.33	180	108772	21.88	ug/L	99
131) hexachlorobutadiene	17.44	225	67180	19.41	ug/L	99
132) naphthalene	17.62	128	195876	22.57	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	95267	22.12	ug/L	96
134) hexachloroethane	15.89	201	60845	19.90	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
Data File : 4B54248.D
Acq On : 5 Aug 2015 12:27 pm
Operator : TOANP
Sample : ic2289-20
Misc : MS88663,V4B2289,w,,,1
ALS Vial : 8 Sample Multiplier: 1

Quant Time: Aug 05 13:44:07 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 13:42:40 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54249.D
 Acq On : 5 Aug 2015 12:54 pm
 Operator : TOANP
 Sample : icc2289-50
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Aug 05 14:13:42 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:12:09 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	142044	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	427186	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	505017	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	457920	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	269700	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	155259	50.55	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	101.10%
50) 1,2-dichloroethane-d4 (s)	9.02	65	169612	49.53	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	99.06%
80) toluene-d8 (s)	11.02	98	598236	50.31	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	100.62%
105) 4-bromofluorobenzene (s)	13.86	95	230299	48.47	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.94%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.67	59	80079	227.25	ug/L
3) 1,4-dioxane	10.11	88	40249	1411.24	ug/L
7) chlorodifluoromethane	3.64	51	204610	49.35	ug/L
8) dichlorodifluoromethane	3.62	85	202349	55.18	ug/L
10) chloromethane	3.96	52	75060	47.70	ug/L
11) vinyl chloride	4.17	62	225178	52.52	ug/L
12) bromomethane	4.75	94	128108	47.69	ug/L
13) chloroethane	4.91	64	89843	50.09	ug/L
15) trichlorofluoromethane	5.28	101	217084	55.10	ug/L
19) ethyl ether	5.63	74	80145	49.88	ug/L
20) 2-chloropropane	5.81	39	23988	37.96	ug/L
21) acrolein	5.86	56	285912	548.23	ug/L
22) 1,1-dichloroethene	6.02	96	156381	45.81	ug/L
23) acetone	6.05	58	14461	52.14	ug/L
24) allyl chloride	6.46	76	115265	47.86	ug/L
25) acetonitrile	6.46	40	132992	485.39	ug/L
26) iodomethane	6.27	142	313347	49.75	ug/L
27) carbon disulfide	6.38	76	472706	47.68	ug/L
28) methylene chloride	6.64	84	176519	49.76	ug/L
29) methyl acetate	6.44	74	25462	56.16	ug/L
30) 1-chloropropane	6.66	42	278651	41.14	ug/L
31) methyl tert butyl ether	6.89	73	441852	51.99	ug/L
32) trans-1,2-dichloroethene	6.95	96	147473	43.05	ug/L
33) di-isopropyl ether	7.41	45	584015	45.46	ug/L
34) 2-butanone	8.08	72	18580	57.13	ug/L
35) 1,1-dichloroethane	7.45	63	291836	50.92	ug/L
36) chloroprene	7.55	53	213013	48.22	ug/L
37) acrylonitrile	6.92	53	316441	296.77	ug/L
38) vinyl acetate	7.43	86	23849	50.31	ug/L
39) ethyl tert-butyl ether	7.83	59	524283	45.68	ug/L
40) ethyl acetate	8.09	45	25157	49.32	ug/L
41) 2,2-dichloropropane	8.11	77	115985	41.45	ug/L
42) cis-1,2-dichloroethene	8.11	96	170444	52.70	ug/L

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54249.D
 Acq On : 5 Aug 2015 12:54 pm
 Operator : TOANP
 Sample : icc2289-50
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Aug 05 14:13:42 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:12:09 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.17	85	24493	55.11	ug/L	97
44) propionitrile	8.18	54	252936	501.11	ug/L	89
45) bromochloromethane	8.40	128	91577	53.62	ug/L	100
46) tetrahydrofuran	8.43	42	57255	45.15	ug/L	95
47) chloroform	8.44	85	181475	48.44	ug/L	99
48) T-BUTYL FORMATE	8.46	59	126667	56.88	ug/L	99
51) freon 113	5.96	151	96902	46.27	ug/L	97
52) methacrylonitrile	8.34	41	109367	48.47	ug/L	98
53) 1,1,1-trichloroethane	8.67	97	185743	49.00	ug/L	99
54) cyclohexane	8.74	84	196417	48.17	ug/L	98
55) iso-butyl alcohol	8.83	43	64292	412.83	ug/L	98
57) epichlorohydrin	10.63	57	86302	252.74	ug/L	99
58) n-butyl alcohol	9.53	56	223927	2257.95	ug/L	99
59) carbon tetrachloride	8.86	117	187904	48.72	ug/L	100
60) 1,1-dichloropropene	8.84	75	203280	49.90	ug/L	100
61) hexane	7.20	57	150307	45.65	ug/L	99
63) benzene	9.08	78	611437	45.34	ug/L	100
64) iso-octane	9.07	57	484301	46.40	ug/L	99
65) tert-amyl methyl ether	9.09	87	103480	46.63	ug/L	95
66) heptane	9.21	57	105867	47.49	ug/L	98
67) isopropyl acetate	8.99	61	69968	20.29	ug/L	96
68) 1,2-dichloroethane	9.10	62	213481	51.32	ug/L	100
69) trichloroethene	9.74	95	167725	51.31	ug/L	98
71) ethyl acrylate	9.73	55	241438	51.43	ug/L	99
72) 2-nitropropane	10.50	41	68646	47.13	ug/L	96
73) 2-chloroethyl vinyl ether	10.50	63	542897	261.29	ug/L	100
74) methyl methacrylate	9.99	100	42697	51.84	ug/L	# 83
75) 1,2-dichloropropane	10.00	63	182101	49.40	ug/L	99
76) dibromomethane	10.16	93	111215	50.94	ug/L	96
77) methylcyclohexane	9.94	83	210115	49.10	ug/L	97
78) bromodichloromethane	10.28	83	238567	53.44	ug/L	99
79) cis-1,3-dichloropropene	10.72	75	298139	52.00	ug/L	99
81) 4-methyl-2-pentanone	10.82	58	71442	51.82	ug/L	98
82) toluene	11.09	92	389476	49.90	ug/L	99
83) 3-methyl-1-butanol	10.83	55	146795	775.95	ug/L	93
84) trans-1,3-dichloropropene	11.29	75	273734	54.05	ug/L	97
85) ethyl methacrylate	11.27	69	226985	58.15	ug/L	99
86) 1,1,2-trichloroethane	11.52	83	142996	51.28	ug/L	99
87) 2-hexanone	11.69	58	64172	53.28	ug/L	91
89) tetrachloroethene	11.69	164	179864	46.18	ug/L	99
90) 1,3-dichloropropane	11.71	76	263602	51.65	ug/L	97
91) butyl acetate	11.76	56	118066	44.07	ug/L	97
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	162954	366.64	ug/L	99
93) dibromochloromethane	11.99	129	203519	51.24	ug/L	100
94) 1,2-dibromoethane	12.15	107	175035	53.47	ug/L	99
96) chlorobenzene	12.63	112	458187	49.48	ug/L	98
97) 1,1,1,2-tetrachloroethane	12.70	131	176304	52.45	ug/L	99
98) ethylbenzene	12.69	91	762305	49.20	ug/L	99
99) m,p-xylene	12.80	106	584841	100.75	ug/L	97
100) o-xylene	13.26	106	304539	52.17	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54249.D
 Acq On : 5 Aug 2015 12:54 pm
 Operator : TOANP
 Sample : icc2289-50
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 9 Sample Multiplier: 1

Quant Time: Aug 05 14:13:42 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:12:09 2015
 Response via : Initial Calibration

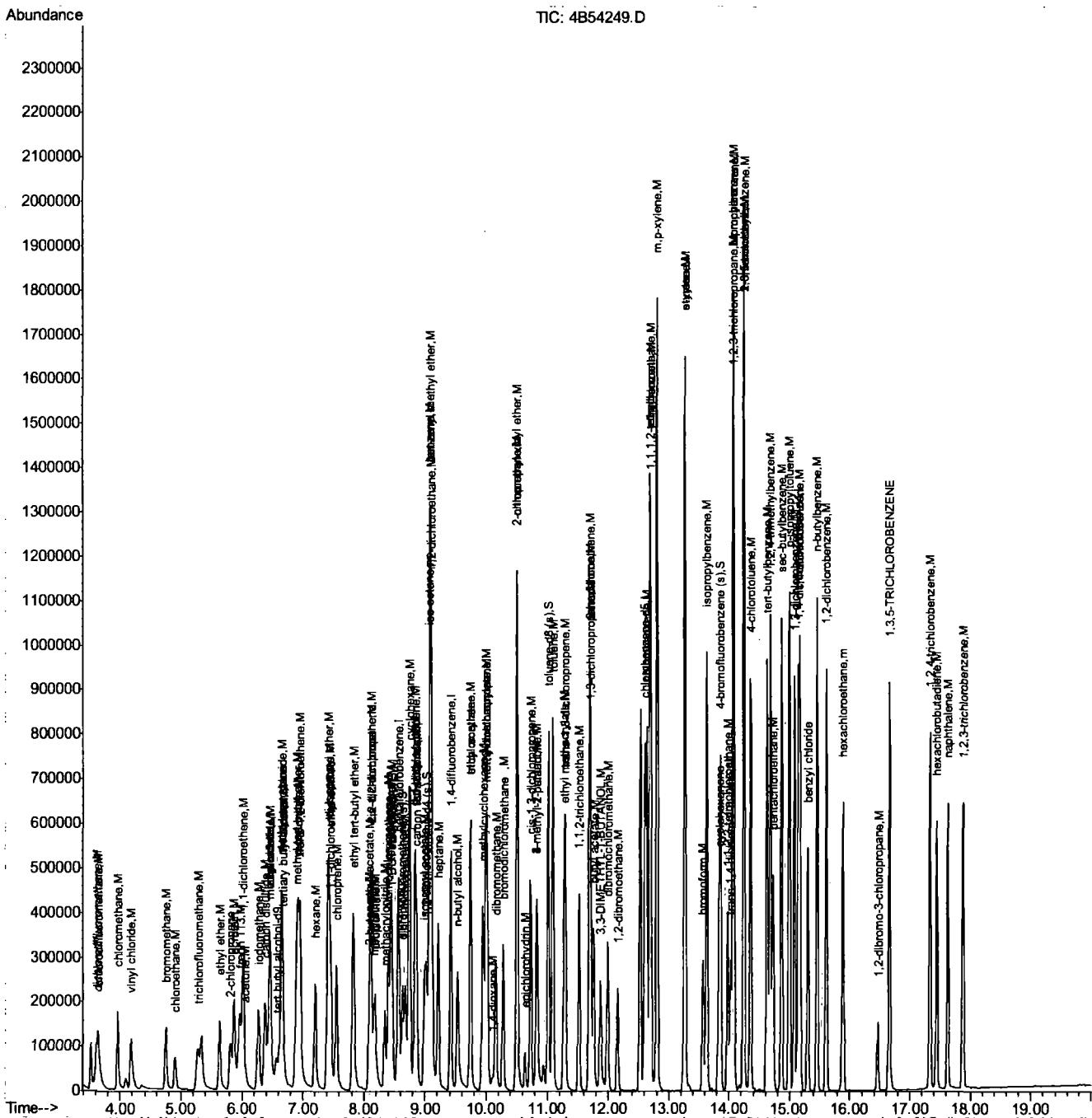
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
101) styrene	13.27	104	514613	54.21	ug/L	98
102) bromoform	13.57	173	162530	55.88	ug/L	98
104) isopropylbenzene	13.63	105	768394	48.19	ug/L	100
106) cyclohexanone	13.83	55	209172	346.13	ug/L	99
107) bromobenzene	14.07	156	235895	48.59	ug/L	95
108) 1,1,2,2-tetrachloroethane	13.98	83	240168	50.08	ug/L	99
109) trans-1,4-dichloro-2-buten	14.02	53	62608	49.02	ug/L	98
110) 1,2,3-trichloropropane	14.06	110	57598	48.25	ug/L	96
111) n-propylbenzene	14.08	91	878004	46.90	ug/L	99
113) 2-chlorotoluene	14.24	126	198745	49.47	ug/L	97
114) 4-chlorotoluene	14.35	91	583335	47.55	ug/L	100
115) 1,3,5-trimethylbenzene	14.24	105	655685	48.49	ug/L	99
116) tert-butylbenzene	14.62	119	578217	47.75	ug/L	98
117) pentachloroethane	14.72	167	120723	46.31	ug/L	97
118) 1,2,4-trimethylbenzene	14.68	105	658527	48.92	ug/L	99
119) sec-butylbenzene	14.87	105	869543	49.43	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	442128	48.44	ug/L	99
121) p-isopropyltoluene	15.00	119	736660	49.85	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	436957	47.69	ug/L	99
123) benzyl chloride	15.30	91	440384	42.39	ug/L	100
124) 1,2-dichlorobenzene	15.61	146	433270	49.83	ug/L	100
126) n-butylbenzene	15.46	92	374005	52.85	ug/L	98
128) 1,2-dibromo-3-chloropropan	16.47	75	37702	46.46	ug/L	93
129) 1,3,5-TRICHLOROBENZENE	16.65	180	352015	51.88	ug/L	99
130) 1,2,4-trichlorobenzene	17.33	180	288084	55.62	ug/L	99
131) hexachlorobutadiene	17.44	225	156639	44.21	ug/L	98
132) naphthalene	17.62	128	516010	56.79	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	241392	53.71	ug/L	99
134) hexachloroethane	15.89	201	151665	48.29	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54249.D
Acq On : 5 Aug 2015 12:54 pm
Operator : TOANP
Sample : iccc2289-50
Misc : MS88663,V4B2289,w,,,1
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Aug 05 14:13:42 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 14:12:09 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54250.D
 Acq On : 5 Aug 2015 1:22 pm
 Operator : TOANP
 Sample : ic2289-100
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Aug 05 14:09:33 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:07:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	152324	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	440342	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	524938	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	475546	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	278457	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	162090	51.40	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	102.80%
50) 1,2-dichloroethane-d4 (s)	9.02	65	172177	49.20	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	98.40%
80) toluene-d8 (s)	11.02	98	618173	50.01	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	100.02%
105) 4-bromofluorobenzene (s)	13.86	95	237181	48.22	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.44%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.69	59	183377	488.82	ug/L # 58
3) 1,4-dioxane	10.11	88	93245	3201.99	ug/L 97
7) chlorodifluoromethane	3.64	51	444171	105.30	ug/L 99
8) dichlorodifluoromethane	3.61	85	444112	119.20	ug/L 100
10) chloromethane	3.96	52	169765	103.54	ug/L 99
11) vinyl chloride	4.18	62	503146	114.47	ug/L 97
12) bromomethane	4.74	94	270335	87.48	ug/L 98
13) chloroethane	4.90	64	194231	104.70	ug/L 98
15) trichlorofluoromethane	5.28	101	482379	120.37	ug/L 97
19) ethyl ether	5.63	74	174510	105.32	ug/L 100
20) 2-chloropropane	5.81	39	48196	71.34	ug/L 80
21) acrolein	5.86	56	624684	1166.98	ug/L 99
22) 1,1-dichloroethene	6.02	96	340104	95.70	ug/L 99
23) acetone	6.06	58	30926	108.91	ug/L 99
24) allyl chloride	6.46	76	274071	108.05	ug/L 91
25) acetonitrile	6.45	40	280810	1010.23	ug/L 94
26) iodomethane	6.27	142	676656	104.00	ug/L 98
27) carbon disulfide	6.38	76	1031150	100.85	ug/L 98
28) methylene chloride	6.64	84	374677	102.08	ug/L 98
29) methyl acetate	6.44	74	54327	119.73	ug/L 96
30) 1-chloropropane	6.66	42	599198	83.21	ug/L 97
31) methyl tert butyl ether	6.89	73	946343	108.89	ug/L 98
32) trans-1,2-dichloroethene	6.95	96	324123	89.89	ug/L 95
33) di-isopropyl ether	7.41	45	1227486	91.36	ug/L 83
34) 2-butanone	8.08	72	42130	132.98	ug/L # 86
35) 1,1-dichloroethane	7.45	63	627638	107.49	ug/L 99
36) chloroprene	7.55	53	467988	103.04	ug/L 99
37) acrylonitrile	6.92	53	668696	626.09	ug/L 98
38) vinyl acetate	7.43	86	51509	106.21	ug/L 78
39) ethyl tert-butyl ether	7.83	59	1108286	92.23	ug/L 99
40) ethyl acetate	8.08	45	53614	101.24	ug/L # 88
41) 2,2-dichloropropane	8.11	77	239671	81.82	ug/L 95
42) cis-1,2-dichloroethene	8.11	96	368672	111.60	ug/L 95

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54250.D
 Acq On : 5 Aug 2015 1:22 pm
 Operator : TOANP
 Sample : ic2289-100
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Aug 05 14:09:33 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:07:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
43) methylacrylate	8.16	85	53820	120.55	ug/L	97
44) propionitrile	8.19	54	552171	1064.70	ug/L	97
45) bromochloromethane	8.40	128	197037	113.42	ug/L	98
46) tetrahydrofuran	8.43	42	123798	93.67	ug/L	98
47) chloroform	8.44	85	379546	97.77	ug/L	99
48) T-BUTYL FORMATE	8.46	59	270320	119.19	ug/L	99
51) freon 113	5.96	151	204380	93.83	ug/L	99
52) methacrylonitrile	8.34	41	240185	102.89	ug/L	96
53) 1,1,1-trichloroethane	8.67	97	402495	103.55	ug/L	99
54) cyclohexane	8.74	84	440505	105.00	ug/L	96
55) iso-butyl alcohol	8.83	43	150898	914.33	ug/L	98
57) epichlorohydrin	10.64	57	188130	489.98	ug/L	98
58) n-butyl alcohol	9.53	56	519748	4005.05	ug/L	99
59) carbon tetrachloride	8.86	117	414271	103.74	ug/L	100
60) 1,1-dichloropropene	8.83	75	451545	107.97	ug/L	99
61) hexane	7.20	57	336409	97.76	ug/L	99
63) benzene	9.08	78	1326122	93.61	ug/L	99
64) iso-octane	9.07	57	1059788	104.26	ug/L	99
65) tert-amyl methyl ether	9.09	87	206475	88.50	ug/L	93
66) heptane	9.21	57	236489	102.47	ug/L	97
67) isopropyl acetate	8.99	61	146886	36.79	ug/L	97
68) 1,2-dichloroethane	9.10	62	440523	102.84	ug/L	99
69) trichloroethene	9.74	95	364370	107.81	ug/L	98
71) ethyl acrylate	9.73	55	519917	106.25	ug/L	99
72) 2-nitropropane	10.50	41	145742	95.70	ug/L	92
73) 2-chloroethyl vinyl ether	10.50	63	1145967	532.84	ug/L	99
74) methyl methacrylate	9.99	100	92959	109.32	ug/L	# 84
75) 1,2-dichloropropane	10.00	63	384852	100.73	ug/L	100
76) dibromomethane	10.16	93	237351	104.67	ug/L	99
77) methylcyclohexane	9.94	83	478713	108.38	ug/L	97
78) bromodichloromethane	10.28	83	509811	110.83	ug/L	99
79) cis-1,3-dichloropropene	10.73	75	645764	109.10	ug/L	98
81) 4-methyl-2-pentanone	10.82	58	155389	110.17	ug/L	98
82) toluene	11.09	92	856288	105.80	ug/L	99
83) 3-methyl-1-butanol	10.83	55	327046	1619.54	ug/L	99
84) trans-1,3-dichloropropene	11.29	75	584815	112.41	ug/L	97
85) ethyl methacrylate	11.27	69	500966	125.70	ug/L	98
86) 1,1,2-trichloroethane	11.52	83	308391	106.24	ug/L	98
87) 2-hexanone	11.69	58	148631	120.51	ug/L	98
89) tetrachloroethene	11.69	164	396281	97.64	ug/L	98
90) 1,3-dichloropropane	11.71	76	558281	106.09	ug/L	98
91) butyl acetate	11.76	56	257559	90.86	ug/L	94
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	364614	766.78	ug/L	98
93) dibromochloromethane	11.99	129	444726	108.08	ug/L	100
94) 1,2-dibromoethane	12.15	107	383789	114.03	ug/L	99
96) chlorobenzene	12.64	112	1008232	105.13	ug/L	100
97) 1,1,1,2-tetrachloroethane	12.70	131	376160	108.95	ug/L	98
98) ethylbenzene	12.69	91	1665552	103.81	ug/L	99
99) m,p-xylene	12.80	106	1296761	216.27	ug/L	96
100) o-xylene	13.26	106	674677	112.17	ug/L	96

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54250.D
 Acq On : 5 Aug 2015 1:22 pm
 Operator : TOANP
 Sample : ic2289-100
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Aug 05 14:09:33 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:07:53 2015
 Response via : Initial Calibration

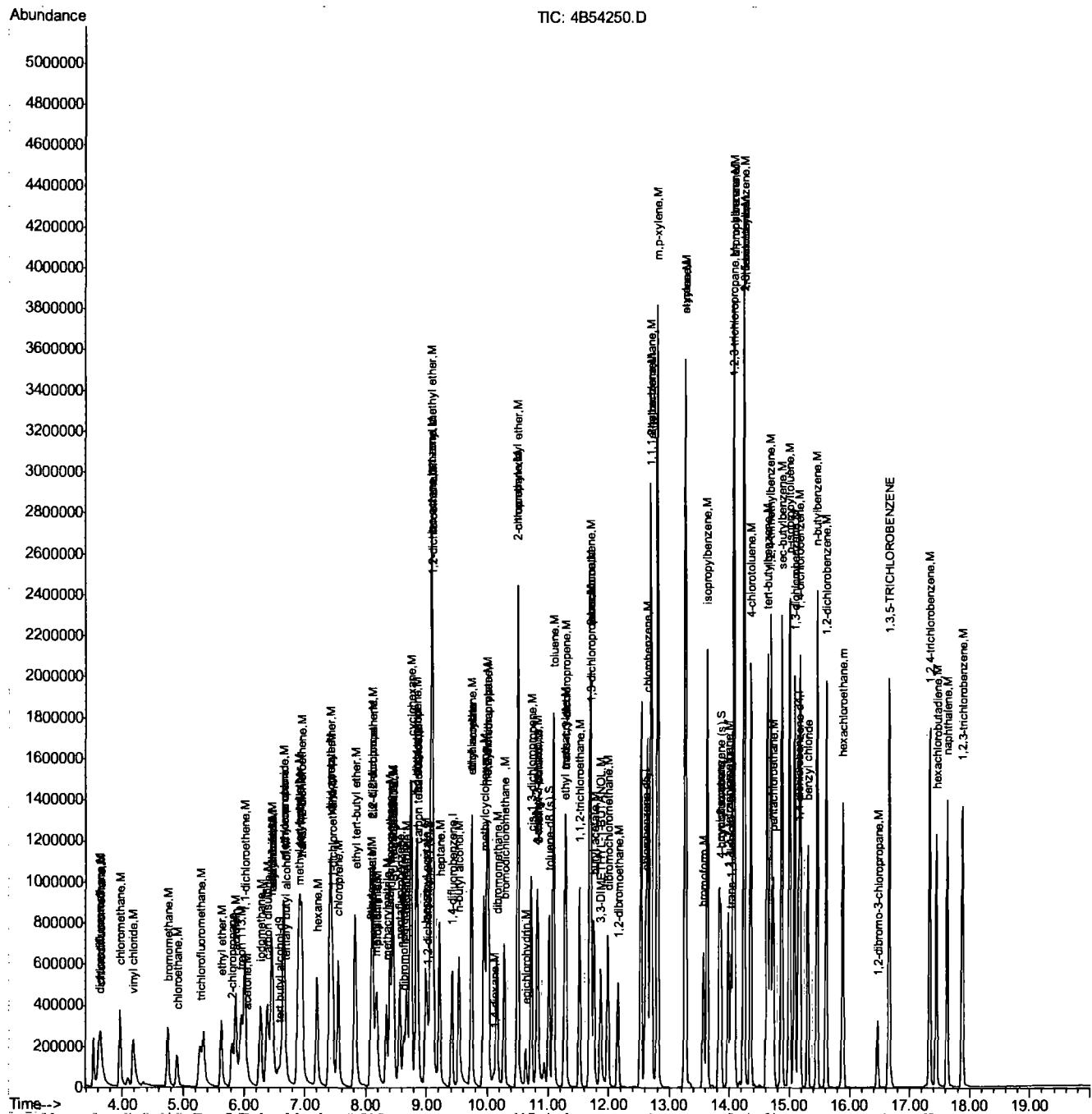
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
101) styrene	13.27	104	1140164	117.34	ug/L	98
102) bromoform	13.57	173	358118	120.26	ug/L	100
104) isopropylbenzene	13.63	105	1686948	102.61	ug/L	99
106) cyclohexanone	13.83	55	514116	800.37	ug/L	98
107) bromobenzene	14.07	156	515647	102.81	ug/L	97
108) 1,1,2,2-tetrachloroethane	13.98	83	515882	104.52	ug/L	99
109) trans-1,2-dichloro-2-buten	14.02	53	137862	104.39	ug/L	99
110) 1,2,3-trichloropropane	14.06	110	122399	99.93	ug/L	97
111) n-propylbenzene	14.08	91	1945144	100.42	ug/L	100
113) 2-chlorotoluene	14.24	126	439544	106.54	ug/L	95
114) 4-chlorotoluene	14.35	91	1270141	99.87	ug/L	100
115) 1,3,5-trimethylbenzene	14.24	105	1437181	103.06	ug/L	100
116) tert-butylbenzene	14.63	119	1258947	100.45	ug/L	96
117) pentachloroethane	14.72	167	257405	94.53	ug/L	98
118) 1,2,4-trimethylbenzene	14.68	105	1412075	101.70	ug/L	100
119) sec-butylbenzene	14.87	105	1902828	105.19	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	950076	100.49	ug/L	99
121) p-isopropyltoluene	15.00	119	1607909	105.97	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	943626	99.13	ug/L	99
123) benzyl chloride	15.30	91	931217	84.47	ug/L	100
124) 1,2-dichlorobenzene	15.61	146	927595	103.54	ug/L	99
126) n-butylbenzene	15.46	92	825537	114.67	ug/L	97
128) 1,2-dibromo-3-chloropropan	16.46	75	83271	98.47	ug/L	90
129) 1,3,5-TRICHLOROBENZENE	16.65	180	747739	107.59	ug/L	100
130) 1,2,4-trichlorobenzene	17.33	180	622234	117.93	ug/L	100
131) hexachlorobutadiene	17.44	225	320403	87.22	ug/L	97
132) naphthalene	17.62	128	1142135	123.99	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	519343	113.62	ug/L	97
134) hexachloroethane	15.89	201	333252	102.68	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54250.D
Acq On : 5 Aug 2015 1:22 pm
Operator : TOANP
Sample : ic2289-100
Misc : MS88663,V4B2289,w,,,1
ALS Vial : 10 Sample Multiplier: 1

Quant Time: Aug 05 14:09:33 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 14:07:53 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54251.D
 Acq On : 5 Aug 2015 1:50 pm
 Operator : TOANP
 Sample : ic2289-200
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Aug 05 14:17:08 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:15:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.59	65	161277	500.00	ug/L	0.02
5) pentafluorobenzene	8.56	168	446123	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	534451	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	480789	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	280134	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	163959	50.92	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	101.84%
50) 1,2-dichloroethane-d4 (s)	9.02	65	169371	47.54	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	95.08%
80) toluene-d8 (s)	11.02	98	624262	49.57	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	99.14%
105) 4-bromofluorobenzene (s)	13.86	95	240197	49.02	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.04%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.69	59	368528	937.24	ug/L # 53
3) 1,4-dioxane	10.11	88	191696	5639.08	ug/L 96
7) chlorodifluoromethane	3.64	51	901529	207.52	ug/L 99
8) dichlorodifluoromethane	3.61	85	850435	214.60	ug/L 97
10) chloromethane	3.96	52	344977	209.92	ug/L 96
11) vinyl chloride	4.18	62	985164	215.52	ug/L 98
12) bromomethane	4.74	94	451750	162.44	ug/L 97
13) chloroethane	4.90	64	343959	182.43	ug/L 95
15) trichlorofluoromethane	5.27	101	916972	215.08	ug/L 97
19) ethyl ether	5.63	74	357920	211.93	ug/L 94
20) 2-chloropropane	5.81	39	84550	139.78	ug/L 87
21) acrolein	5.87	56	1250055	2223.36	ug/L 98
22) 1,1-dichloroethene	6.02	96	720334	205.07	ug/L 93
23) acetone	6.06	58	60091	203.25	ug/L # 81
24) allyl chloride	6.46	76	573944	226.20	ug/L # 81
25) acetonitrile	6.46	40	558194	1962.22	ug/L 99
26) iodomethane	6.27	142	1400146	211.88	ug/L 98
27) carbon disulfide	6.38	76	2117922	205.50	ug/L 99
28) methylene chloride	6.64	84	774609	208.59	ug/L 95
29) methyl acetate	6.44	74	113584	230.48	ug/L 90
30) 1-chloropropane	6.66	42	1193835	175.78	ug/L 99
31) methyl tert butyl ether	6.89	73	1907117	212.03	ug/L 99
32) trans-1,2-dichloroethene	6.95	96	670424	192.72	ug/L 97
33) di-isopropyl ether	7.41	45	2414584	183.30	ug/L 88
34) 2-butanone	8.08	72	83894	231.59	ug/L # 87
35) 1,1-dichloroethane	7.45	63	1253371	207.54	ug/L 99
36) chloroprene	7.54	53	928786	201.52	ug/L 98
37) acrylonitrile	6.93	53	1365697	1173.75	ug/L 98
38) vinyl acetate	7.43	86	105799	211.90	ug/L 95
39) ethyl tert-butyl ether	7.83	59	2228734	189.10	ug/L 100
40) ethyl acetate	8.09	45	105610	198.09	ug/L 87
41) 2,2-dichloropropane	8.11	77	435580	155.69	ug/L 94
42) cis-1,2-dichloroethene	8.11	96	751583	218.65	ug/L 94

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54251.D
 Acq On : 5 Aug 2015 1:50 pm
 Operator : TOANP
 Sample : ic2289-200
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Aug 05 14:17:08 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:15:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.17	85	111432	229.48	ug/L	95
44) propionitrile	8.19	54	1125606	2118.56	ug/L	92
45) bromochloromethane	8.40	128	407776	223.29	ug/L	95
46) tetrahydrofuran	8.43	42	241984	186.73	ug/L	99
47) chloroform	8.44	85	785999	201.97	ug/L	96
48) T-BUTYL FORMATE	8.47	59	555520	230.79	ug/L	97
51) freon 113	5.96	151	415553	193.53	ug/L	98
52) methacrylonitrile	8.34	41	480235	203.75	ug/L	98
53) 1,1,1-trichloroethane	8.67	97	799081	201.61	ug/L	99
54) cyclohexane	8.74	84	919328	215.59	ug/L	90
55) iso-butyl alcohol	8.84	43	295848	1873.94	ug/L	98
57) epichlorohydrin	10.64	57	385542	1057.50	ug/L	99
58) n-butyl alcohol	9.53	56	1060475	10217.23	ug/L	98
59) carbon tetrachloride	8.86	117	847352	207.39	ug/L	98
60) 1,1-dichloropropene	8.84	75	917331	211.27	ug/L	98
61) hexane	7.20	57	668510	194.37	ug/L	98
63) benzene	9.08	78	2704139	192.62	ug/L	99
64) iso-octane	9.07	57	2126092	194.81	ug/L	95
65) tert-amyl methyl ether	9.09	87	436196	189.37	ug/L	95
66) heptane	9.21	57	472770	201.14	ug/L	99
67) isopropyl acetate	8.99	61	296166	95.26	ug/L	100
68) 1,2-dichloroethane	9.10	62	864345	195.38	ug/L	98
69) trichloroethene	9.74	95	743930	212.71	ug/L	98
71) ethyl acrylate	9.73	55	1020839	203.34	ug/L	98
72) 2-nitropropane	10.50	41	279130	183.27	ug/L	92
73) 2-chloroethyl vinyl ether	10.50	63	2272973	1021.62	ug/L	97
74) methyl methacrylate	9.99	100	193525	218.20	ug/L	94
75) 1,2-dichloropropane	10.00	63	769783	197.52	ug/L	99
76) dibromomethane	10.16	93	484947	208.21	ug/L	98
77) methylcyclohexane	9.94	83	962349	210.96	ug/L	98
78) bromodichloromethane	10.28	83	1040096	216.14	ug/L	99
79) cis-1,3-dichloropropene	10.73	75	1308146	212.69	ug/L	97
81) 4-methyl-2-pentanone	10.82	58	315959	213.34	ug/L	96
82) toluene	11.09	92	1779838	214.20	ug/L	97
83) 3-methyl-1-butanol	10.83	55	654036	3415.76	ug/L	96
84) trans-1,3-dichloropropene	11.29	75	1178558	215.29	ug/L	99
85) ethyl methacrylate	11.28	69	1020068	236.49	ug/L	96
86) 1,1,2-trichloroethane	11.52	83	626099	209.81	ug/L	98
87) 2-hexanone	11.69	58	295973	225.08	ug/L	95
89) tetrachloroethene	11.69	164	795453	196.89	ug/L	99
90) 1,3-dichloropropane	11.71	76	1120916	207.21	ug/L	98
91) butyl acetate	11.76	56	526564	191.31	ug/L	97
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	730469	1652.93	ug/L	97
93) dibromochloromethane	11.99	129	921933	218.27	ug/L	99
94) 1,2-dibromoethane	12.15	107	793606	225.91	ug/L	99
96) chlorobenzene	12.64	112	2088368	213.89	ug/L	98
97) 1,1,1,2-tetrachloroethane	12.70	131	785333	219.45	ug/L	99
98) ethylbenzene	12.69	91	3409958	209.17	ug/L	98
99) m,p-xylene	12.80	106	2688195	437.04	ug/L	96
100) o-xylene	13.26	106	1392519	223.33	ug/L	96

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54251.D
 Acq On : 5 Aug 2015 1:50 pm
 Operator : TOANP
 Sample : ic2289-200
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Aug 05 14:17:08 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Wed Aug 05 14:15:54 2015
 Response via : Initial Calibration

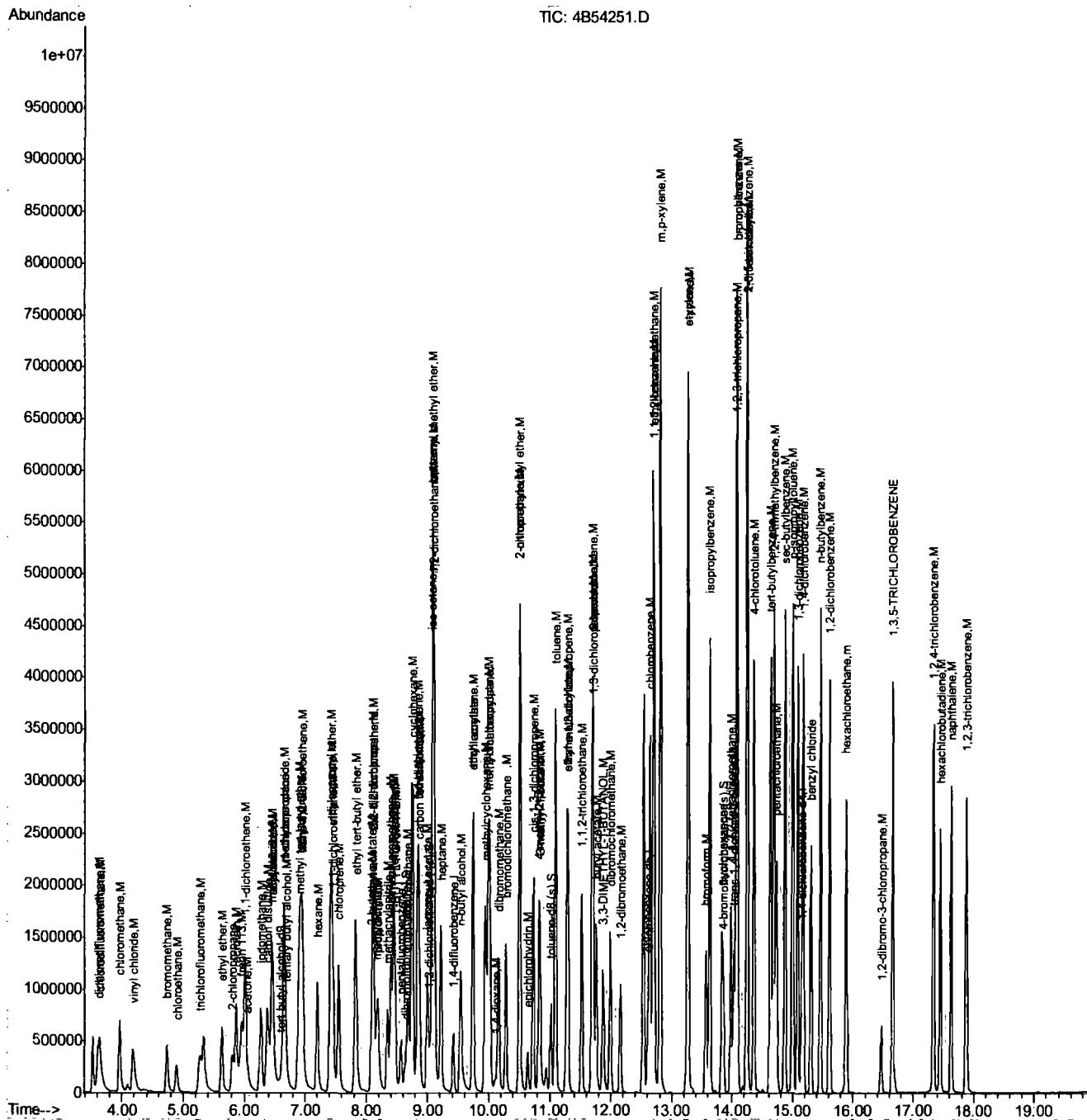
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
101) styrene	13.27	104	2357594	230.36	ug/L	97
102) bromoform	13.57	173	745935	236.30	ug/L	99
104) isopropylbenzene	13.63	105	3487739	210.86	ug/L	98
106) cyclohexanone	13.83	55	817403	1376.21	ug/L	99
107) bromobenzene	14.08	156	1078873	213.92	ug/L	89
108) 1,1,2,2-tetrachloroethane	13.98	83	1058293	211.44	ug/L	99
109) trans-1,4-dichloro-2-buten	14.02	53	280808	210.99	ug/L	97
110) 1,2,3-trichloropropane	14.06	110	247600	200.74	ug/L	97
111) n-propylbenzene	14.08	91	3994495	206.71	ug/L	97
113) 2-chlorotoluene	14.24	126	914110	217.86	ug/L	94
114) 4-chlorotoluene	14.35	91	2630615	207.52	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	2932524	208.82	ug/L	100
116) tert-butylbenzene	14.63	119	2566897	204.97	ug/L	98
117) pentachloroethane	14.72	167	568321	213.00	ug/L	96
118) 1,2,4-trimethylbenzene	14.68	105	2903953	207.81	ug/L	99
119) sec-butylbenzene	14.87	105	3910378	213.14	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	1969144	208.24	ug/L	99
121) p-isopropyltoluene	15.00	119	3266745	211.63	ug/L	98
122) 1,4-dichlorobenzene	15.17	146	1956832	206.73	ug/L	98
123) benzyl chloride	15.30	91	1889010	180.78	ug/L	99
124) 1,2-dichlorobenzene	15.61	146	1883480	207.86	ug/L	99
126) n-butylbenzene	15.46	92	1662661	221.58	ug/L	99
128) 1,2-dibromo-3-chloropropan	16.46	75	170232	204.20	ug/L	99
129) 1,3,5-TRICHLOROBENZENE	16.65	180	1514507	212.41	ug/L	98
130) 1,2,4-trichlorobenzene	17.33	180	1292871	233.18	ug/L	99
131) hexachlorobutadiene	17.44	225	666903	186.18	ug/L	98
132) naphthalene	17.62	128	2426040	247.36	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	1088592	228.29	ug/L	98
134) hexachloroethane	15.89	201	697707	214.04	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
Data File : 4B54251.D
Acq On : 5 Aug 2015 1:50 pm
Operator : TOANP
Sample : ic2289-200
Misc : MS88663,V4B2289,w,,,1
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Aug 05 14:17:08 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Wed Aug 05 14:15:54 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54254.D
 Acq On : 5 Aug 2015 3:20 pm
 Operator : TOANP
 Sample : icv2289-50
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Aug 10 08:21:56 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.57	65	142025	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	440912	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	520562	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	462444	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	265659	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	159564	50.05	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	100.10%
50) 1,2-dichloroethane-d4 (s)	9.02	65	167138	47.70	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	95.40%
80) toluene-d8 (s)	11.01	98	608517	49.65	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	99.30%
105) 4-bromofluorobenzene (s)	13.86	95	229251	49.43	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.86%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.67	59	87967	256.05	ug/L
3) 1,4-dioxane	10.11	88	44745	1408.78	ug/L
7) chlorodifluoromethane	3.63	51	197661	44.33	ug/L
8) dichlorodifluoromethane	3.60	85	167005	39.11	ug/L
10) chloromethane	3.96	52	74950	45.89	ug/L
11) vinyl chloride	4.17	62	205845	43.86	ug/L
12) bromomethane	4.76	94	136246	50.63	ug/L
13) chloroethane	4.92	64	106922	57.94	ug/L
15) trichlorofluoromethane	5.27	101	204231	46.07	ug/L
19) ethyl ether	5.63	74	83783	49.87	ug/L
20) 2-chloropropane	5.81	39	17729	33.16	ug/L
21) acrolein	5.86	56	351480	624.78	ug/L
22) 1,1-dichloroethene	6.03	96	170623	49.01	ug/L
23) acetone	6.05	58	15821	54.02	ug/L
24) allyl chloride	6.45	76	162777	63.87	ug/L
25) acetonitrile	6.45	40	137193	489.30	ug/L
26) iodomethane	6.28	142	302743	46.05	ug/L
27) carbon disulfide	6.39	76	476624	46.65	ug/L
28) methylene chloride	6.65	84	185683	50.35	ug/L
29) methyl acetate	6.44	74	24423	49.21	ug/L
30) 1-chloropropane	6.65	42	286588	45.54	ug/L
31) methyl tert butyl ether	6.89	73	878876	94.61	ug/L
32) trans-1,2-dichloroethene	6.95	96	153059	44.70	ug/L
33) di-isopropyl ether	7.41	45	588209	48.28	ug/L
34) 2-butanone	8.08	72	19634	50.31	ug/L
35) 1,1-dichloroethane	7.45	63	304088	47.26	ug/L
36) chloroprene	7.54	53	214783	45.07	ug/L
37) acrylonitrile	6.92	53	333935	254.82	ug/L
38) vinyl acetate	7.43	86	26891	54.09	ug/L
39) ethyl tert-butyl ether	7.83	59	534086	49.35	ug/L
40) ethyl acetate	8.08	45	25023	47.24	ug/L
41) 2,2-dichloropropane	8.10	77	105621	38.82	ug/L
42) cis-1,2-dichloroethene	8.11	96	174961	47.04	ug/L

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54254.D
 Acq On : 5 Aug 2015 3:20 pm
 Operator : TOANP
 Sample : icv2289-50
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Aug 10 08:21:56 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.16	85	25186	49.86	ug/L	99
44) propionitrile	8.18	54	251570	475.95	ug/L	94
45) bromochloromethane	8.40	128	94204	49.35	ug/L	98
46) tetrahydrofuran	8.43	42	59813	47.09	ug/L	98
47) chloroform	8.44	85	188129	48.86	ug/L	96
48) T-BUTYL FORMATE	8.46	59	123560	50.21	ug/L	97
51) freon 113	5.97	151	97918	46.33	ug/L	100
52) methacrylonitrile	8.34	41	114391	49.00	ug/L	99
53) 1,1,1-trichloroethane	8.67	97	184506	45.32	ug/L	99
54) cyclohexane	8.73	84	173386	38.87	ug/L	92
55) iso-butyl alcohol	8.82	43	68583	472.12	ug/L	97
57) epichlorohydrin	10.63	57	90757	253.96	ug/L	98
58) n-butyl alcohol	9.53	56	242937	2397.26	ug/L	100
59) carbon tetrachloride	8.86	117	183882	44.14	ug/L	97
60) 1,1-dichloropropene	8.83	75	218404	47.19	ug/L	100
61) hexane	7.20	57	190194	54.10	ug/L	98
63) benzene	9.08	78	638426	46.86	ug/L	99
64) iso-octane	9.07	57	384768	35.16	ug/L	96
65) tert-amyl methyl ether	9.09	87	100495	46.40	ug/L	96
66) heptane	9.21	57	126451	52.25	ug/L	96
67) isopropyl acetate	8.99	61	70195	43.10	ug/L	92
68) 1,2-dichloroethane	9.10	62	213155	47.49	ug/L	99
69) trichloroethene	9.74	95	172718	47.11	ug/L	98
71) ethyl acrylate	9.73	55	241314	45.80	ug/L	99
72) 2-nitropropane	10.50	41	72108	49.06	ug/L	96
73) 2-chloroethyl vinyl ether	10.50	63	606109	279.09	ug/L	99
74) methyl methacrylate	9.99	100	46233	52.92	ug/L	98
75) 1,2-dichloropropane	10.00	63	188043	48.05	ug/L	99
76) dibromomethane	10.16	93	111139	48.77	ug/L	98
77) methylcyclohexane	9.94	83	206332	43.76	ug/L	99
78) bromodichloromethane	10.28	83	243498	49.69	ug/L	100
79) cis-1,3-dichloropropene	10.72	75	303936	47.87	ug/L	99
81) 4-methyl-2-pentanone	10.82	58	76491	50.85	ug/L	96
82) toluene	11.09	92	404953	46.64	ug/L	99
83) 3-methyl-1-butanol	10.83	55	157924	940.34	ug/L	96
84) trans-1,3-dichloropropene	11.29	75	261048	46.95	ug/L	98
85) ethyl methacrylate	11.27	69	230523	50.11	ug/L	99
86) 1,1,2-trichloroethane	11.52	83	146584	48.50	ug/L	97
87) 2-hexanone	11.69	58	68053	49.55	ug/L	95
89) tetrachloroethene	11.69	164	184446	45.91	ug/L	99
90) 1,3-dichloropropane	11.71	76	264286	48.56	ug/L	99
91) butyl acetate	11.76	56	120141	48.39	ug/L	98
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	160602	440.84	ug/L	100
93) dibromochloromethane	11.99	129	207803	48.82	ug/L	99
94) 1,2-dibromoethane	12.15	107	178097	49.95	ug/L	99
96) chlorobenzene	12.63	112	477070	48.78	ug/L	100
97) 1,1,1,2-tetrachloroethane	12.70	131	174915	46.79	ug/L	98
98) ethylbenzene	12.69	91	768867	47.20	ug/L	99
99) m,p-xylene	12.80	106	598309	96.38	ug/L	98
100) o-xylene	13.25	106	313035	50.02	ug/L	97

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54254.D
 Acq On : 5 Aug 2015 3:20 pm
 Operator : TOANP
 Sample : icv2289-50
 Misc : MS88663,V4B2289,w,,,1
 ALS Vial : 14 Sample Multiplier: 1

Quant Time: Aug 10 08:21:56 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

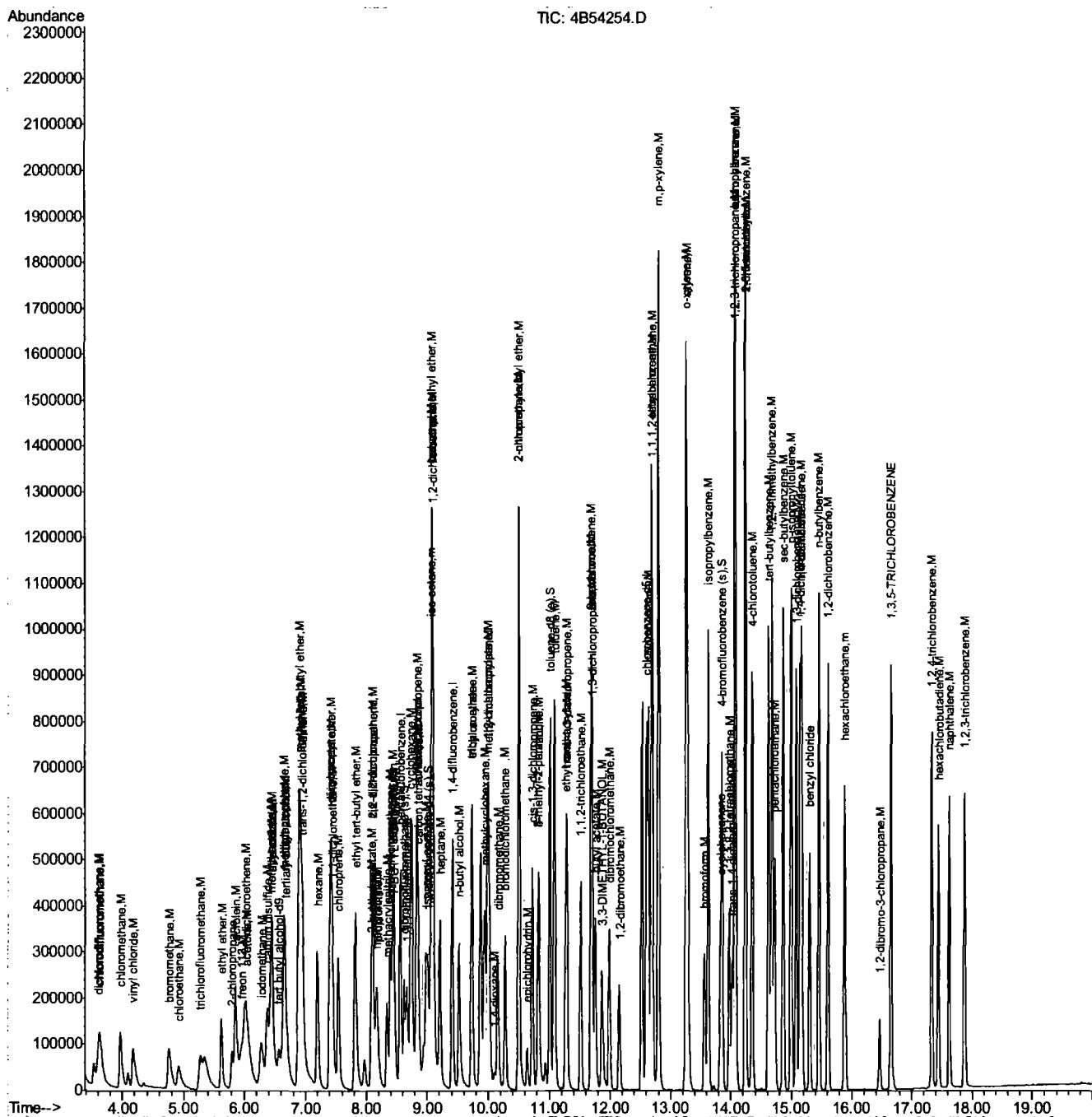
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
101) styrene	13.27	104	522540	48.90	ug/L	98
102) bromoform	13.57	173	165988	50.18	ug/L	98
104) isopropylbenzene	13.63	105	773119	45.84	ug/L	99
106) cyclohexanone	13.83	55	168326	349.81	ug/L	99
107) bromobenzene	14.07	156	235975	47.70	ug/L	99
108) 1,1,2,2-tetrachloroethane	13.98	83	236372	47.86	ug/L	97
109) trans-1,2-dichloro-2-buten	14.02	53	60439	47.60	ug/L	98
110) 1,2,3-trichloropropane	14.06	110	56748	48.49	ug/L	98
111) n-propylbenzene	14.08	91	930222	48.80	ug/L	99
113) 2-chlorotoluene	14.24	126	198113	45.85	ug/L	95
114) 4-chlorotoluene	14.35	91	576327	46.76	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	655414	46.17	ug/L	99
116) tert-butylbenzene	14.62	119	585576	47.55	ug/L	98
117) pentachloroethane	14.72	167	125448	49.22	ug/L	98
118) 1,2,4-trimethylbenzene	14.68	105	683004	49.95	ug/L	99
119) sec-butylbenzene	14.87	105	858564	45.88	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	438486	48.70	ug/L	99
121) p-isopropyltoluene	15.00	119	736359	47.07	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	433803	48.16	ug/L	98
123) benzyl chloride	15.30	91	423232	47.02	ug/L	100
124) 1,2-dichlorobenzene	15.61	146	434616	49.43	ug/L	100
126) n-butylbenzene	15.46	92	371034	47.82	ug/L	98
128) 1,2-dibromo-3-chloropropan	16.46	75	39033	49.24	ug/L	98
129) 1,3,5-TRICHLOROBENZENE	16.65	180	357544	50.99	ug/L	100
130) 1,2,4-trichlorobenzene	17.33	180	277819	49.46	ug/L	99
131) hexachlorobutadiene	17.44	225	150966	44.16	ug/L	98
132) naphthalene	17.62	128	521854	52.85	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	240865	50.84	ug/L	99
134) hexachloroethane	15.89	201	157343	48.51	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54254.D
Acq On : 5 Aug 2015 3:20 pm
Operator : TOANP
Sample : icv2289-50
Misc : MS88663, V4B2289,w,,,1
ALS Vial : 14 Sample Multiplier: 1

Quant Time: Aug 10 08:21:56 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54407.D
 Acq On : 12 Aug 2015 10:09 am
 Operator : TOANP
 Sample : CC2289-20
 Misc : MS89342,V4B2296,w,,,1
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 12 17:18:21 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) tert butyl alcohol-d9	6.58	65	140756	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	435392	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	495542	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	438556	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	244012	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	154824	49.18	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	= 98.36%	
50) 1,2-dichloroethane-d4 (s)	9.02	65	159525	46.11	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	= 92.22%	
80) toluene-d8 (s)	11.02	98	574436	49.24	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	= 98.48%	
105) 4-bromofluorobenzene (s)	13.86	95	208750	49.00	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 98.00%	

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.68	59	32411	95.19	ug/L # 63
3) 1,4-dioxane	10.11	88	18001	571.87	ug/L 89
7) chlorodifluoromethane	3.63	51	78536	17.84	ug/L 97
8) dichlorodifluoromethane	3.61	85	74045	17.56	ug/L 98
10) chloromethane	3.97	52	30047	18.63	ug/L 95
11) vinyl chloride	4.17	62	79691	17.20	ug/L 97
12) bromomethane	4.76	94	55553	20.90	ug/L 96
13) chloroethane	4.92	64	38221	20.98	ug/L 97
15) trichlorofluoromethane	5.28	101	87270	19.94	ug/L 97
19) ethyl ether	5.63	74	28463	17.16	ug/L 92
20) 2-chloropropane	5.81	39	9743	18.45	ug/L 79
21) acrolein	5.86	56	91752	165.16	ug/L 94
22) 1,1-dichloroethene	6.03	96	65076	18.93	ug/L 96
23) acetone	6.05	58	4993	17.26	ug/L # 75
24) allyl chloride	6.46	76	46531	18.49	ug/L 93
25) acetonitrile	6.45	40	46949	169.57	ug/L # 70
26) iodomethane	6.28	142	121113	18.66	ug/L 100
27) carbon disulfide	6.38	76	192311	19.06	ug/L 98
28) methylene chloride	6.65	84	67996	18.67	ug/L 91
29) methyl acetate	6.44	74	8834	18.02	ug/L # 79
30) 1-chloropropane	6.66	42	107359	17.28	ug/L 97
31) methyl tert butyl ether	6.89	73	168557	18.38	ug/L 98
32) trans-1,2-dichloroethene	6.95	96	59183	17.50	ug/L 96
33) di-isopropyl ether	7.41	45	198729	16.52	ug/L 97
34) 2-butanone	8.08	72	6298	16.34	ug/L # 94
35) 1,1-dichloroethane	7.46	63	110612	17.41	ug/L 98
36) chloroprene	7.55	53	84661	17.99	ug/L 99
37) acrylonitrile	6.93	53	109907	84.93	ug/L 99
38) vinyl acetate	7.43	86	8917	18.16	ug/L 84
39) ethyl tert-butyl ether	7.83	59	192034	17.97	ug/L 96
40) ethyl acetate	8.08	45	8595	16.43	ug/L # 86
41) 2,2-dichloropropane	8.11	77	49695	18.50	ug/L 92
42) cis-1,2-dichloroethene	8.11	96	66171	18.02	ug/L 93

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54407.D
 Acq On : 12 Aug 2015 10:09 am
 Operator : TOANP
 Sample : CC2289-20
 Misc : MS89342,V4B2296,w,,,1
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 12 17:18:21 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.16	85	8773	17.59	ug/L #	89
44) propionitrile	8.19	54	87507	167.66	ug/L	89
45) bromochloromethane	8.40	128	35744	18.96	ug/L	87
46) tetrahydrofuran	8.43	42	20463	16.32	ug/L	95
47) chloroform	8.44	85	69826	18.37	ug/L	96
48) T-BUTYL FORMATE	8.46	59	44553	18.33	ug/L	97
51) freon 113	5.97	151	38032	18.22	ug/L	94
52) methacrylonitrile	8.34	41	36616	15.88	ug/L	98
53) 1,1,1-trichloroethane	8.67	97	75033	18.67	ug/L	99
54) cyclohexane	8.74	84	69428	15.76	ug/L #	83
55) iso-butyl alcohol	8.83	43	23688	165.14	ug/L	94
57) epichlorohydrin	10.64	57	29391	86.39	ug/L	100
58) n-butyl alcohol	9.53	56	96335	998.61	ug/L	96
59) carbon tetrachloride	8.86	117	71408	18.01	ug/L	98
60) 1,1-dichloropropene	8.83	75	79539	18.05	ug/L	97
61) hexane	7.20	57	56149	16.78	ug/L	94
63) benzene	9.08	78	239816	18.49	ug/L	98
64) iso-octane	9.07	57	161230	15.48	ug/L	99
65) tert-amyl methyl ether	9.09	87	36924	17.91	ug/L	97
66) heptane	9.21	57	39904	17.32	ug/L	98
67) isopropyl acetate	8.99	61	25790	16.64	ug/L #	86
68) 1,2-dichloroethane	9.10	62	82048	19.20	ug/L	93
69) trichloroethene	9.74	95	67521	19.35	ug/L	99
72) 2-nitropropane	10.50	41	24219	17.31	ug/L #	76
73) 2-chloroethyl vinyl ether	10.50	63	174685	84.50	ug/L	98
74) methyl methacrylate	9.99	100	15646	18.81	ug/L	96
75) 1,2-dichloropropane	10.00	63	66617	17.88	ug/L	98
76) dibromomethane	10.16	93	41401	19.08	ug/L	97
77) methylcyclohexane	9.94	83	78127	17.41	ug/L	96
78) bromodichloromethane	10.28	83	89666	19.22	ug/L	100
79) cis-1,3-dichloropropene	10.72	75	107643	17.81	ug/L	97
81) 4-methyl-2-pentanone	10.82	58	25135	17.55	ug/L	98
82) toluene	11.09	92	153089	18.52	ug/L	98
83) 3-methyl-1-butanol	10.83	55	60866	380.72	ug/L	91
84) trans-1,3-dichloropropene	11.29	75	96087	18.15	ug/L	97
85) ethyl methacrylate	11.27	69	77985	17.81	ug/L	96
86) 1,1,2-trichloroethane	11.52	83	52613	18.29	ug/L	97
87) 2-hexanone	11.69	58	21928	16.77	ug/L	89
89) tetrachloroethene	11.69	164	74355	19.52	ug/L	96
90) 1,3-dichloropropane	11.71	76	95390	18.48	ug/L	97
91) butyl acetate	11.76	56	38462	16.34	ug/L	92
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	70834	205.02	ug/L	99
93) dibromochloromethane	11.99	129	76537	18.96	ug/L	98
94) 1,2-dibromoethane	12.15	107	65690	19.43	ug/L	99
96) chlorobenzene	12.63	112	180201	19.43	ug/L	97
97) 1,1,1,2-tetrachloroethane	12.69	131	68537	19.33	ug/L	99
98) ethylbenzene	12.68	91	297416	19.25	ug/L	100
99) m,p-xylene	12.80	106	228496	38.81	ug/L	96
100) o-xylene	13.26	106	118583	19.98	ug/L	97
101) styrene	13.27	104	192333	18.98	ug/L	98

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54407.D
 Acq On : 12 Aug 2015 10:09 am
 Operator : TOANP
 Sample : CC2289-20
 Misc : MS89342,V4B2296,w,,,1
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 12 17:18:21 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

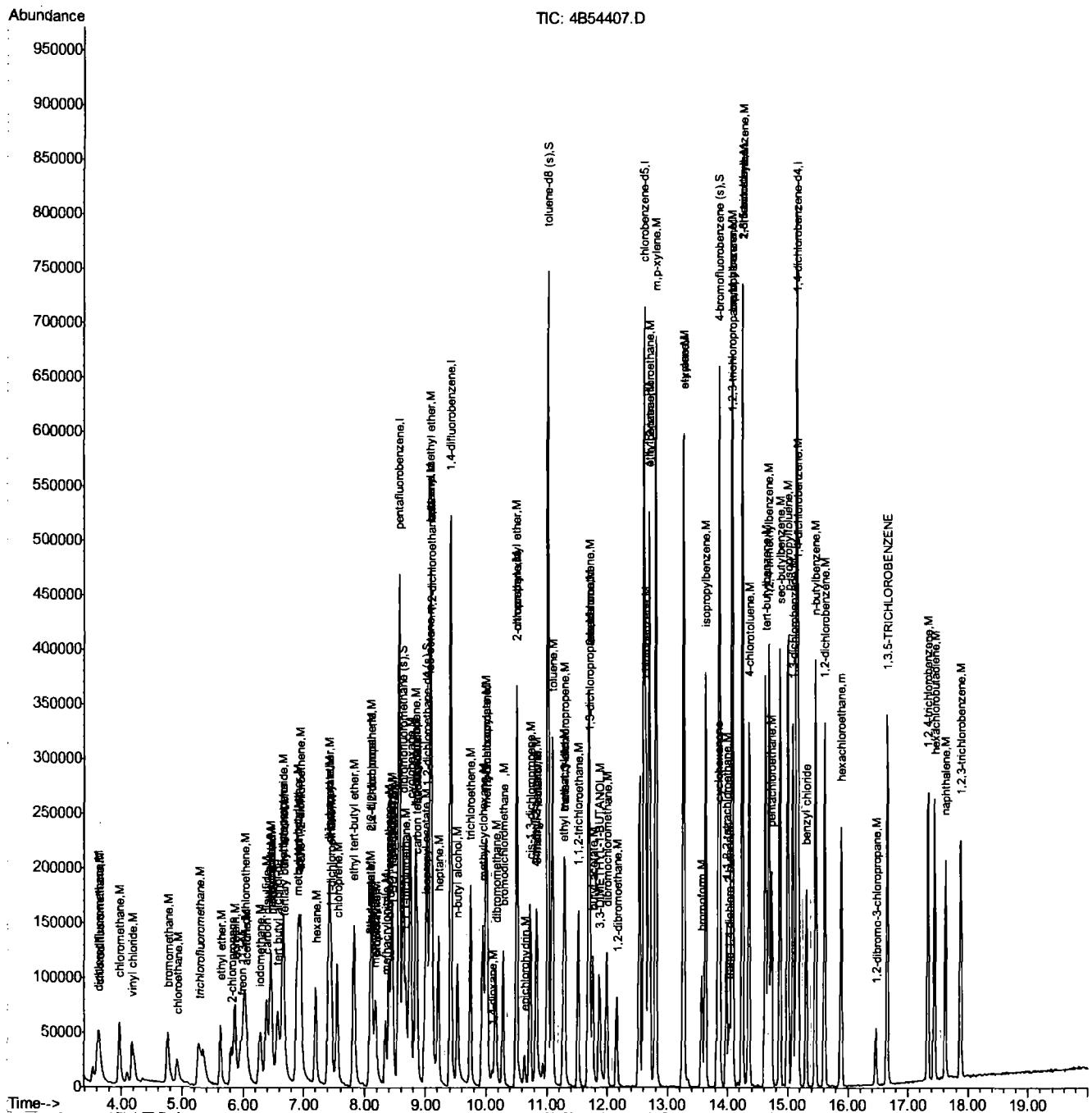
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
102) bromoform	13.57	173	55977	17.84	ug/L	99
104) isopropylbenzene	13.63	105	297794	19.23	ug/L	99
106) cyclohexanone	13.83	55	85294	192.98	ug/L	98
107) bromobenzene	14.07	156	90966	20.02	ug/L	95
108) 1,1,2,2-tetrachloroethane	13.98	83	86876	19.15	ug/L	99
109) trans-1,4-dichloro-2-butene	14.02	53	12970	11.12	ug/L	96
110) 1,2,3-trichloropropane	14.06	110	21298	19.81	ug/L	91
111) n-propylbenzene	14.08	91	337920	19.30	ug/L	99
113) 2-chlorotoluene	14.24	126	76782	19.35	ug/L	97
114) 4-chlorotoluene	14.35	91	213465	18.86	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	253047	19.41	ug/L	96
116) tert-butylbenzene	14.62	119	225930	19.97	ug/L	98
117) pentachloroethane	14.72	167	49325	21.07	ug/L	98
118) 1,2,4-trimethylbenzene	14.68	105	248088	19.75	ug/L	98
119) sec-butylbenzene	14.87	105	329668	19.18	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	163564	19.78	ug/L	99
121) p-isopropyltoluene	15.00	119	283787	19.75	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	161786	19.56	ug/L	99
123) benzyl chloride	15.30	91	146343	17.70	ug/L	99
124) 1,2-dichlorobenzene	15.61	146	162295	20.10	ug/L	98
126) n-butylbenzene	15.46	92	135214	18.97	ug/L	98
128) 1,2-dibromo-3-chloropropan	16.46	75	13367	18.36	ug/L	98
129) 1,3,5-TRICHLOROBENZENE	16.65	180	129430	20.09	ug/L	99
130) 1,2,4-trichlorobenzene	17.33	180	96340	18.67	ug/L	96
131) hexachlorobutadiene	17.44	225	69047	21.99	ug/L	96
132) naphthalene	17.62	128	167117	18.43	ug/L	97
133) 1,2,3-trichlorobenzene	17.88	180	82679	19.00	ug/L	95
134) hexachloroethane	15.89	201	57045	19.15	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54407.D
Acq On : 12 Aug 2015 10:09 am
Operator : TOANP
Sample : CC2289-20
Misc : MS89342,V4B2296,w,,,1
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 12 17:18:21 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54431.D
 Acq On : 12 Aug 2015 9:35 pm
 Operator : TOANP
 Sample : cc2289-50
 Misc : MS89499,V4B2297,w,,,1
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Aug 13 12:17:10 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.58	65	129429	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	415468	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.42	114	477620	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	432693	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	251516	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.63	113	148780	49.52	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	99.04%
50) 1,2-dichloroethane-d4 (s)	9.02	65	159182	48.21	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	96.42%
80) toluene-d8 (s)	11.02	98	556205	49.47	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	98.94%
105) 4-bromofluorobenzene (s)	13.86	95	213058	48.52	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	97.04%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.68	59	84461	269.77	ug/L # 60
3) 1,4-dioxane	10.11	88	39527	1365.61	ug/L 96
7) chlorodifluoromethane	3.64	51	186424	44.37	ug/L 96
8) dichlorodifluoromethane	3.60	85	188678	46.89	ug/L 98
10) chloromethane	3.96	52	65673	42.68	ug/L 94
11) vinyl chloride	4.18	62	192470	43.53	ug/L 98
12) bromomethane	4.75	94	122829	48.44	ug/L 98
13) chloroethane	4.91	64	81645	46.96	ug/L 97
15) trichlorofluoromethane	5.28	101	218002	52.19	ug/L 95
19) ethyl ether	5.62	74	69516	43.91	ug/L 87
20) 2-chloropropane	5.81	39	26758	53.11	ug/L 76
21) acrolein	5.86	56	251766	474.94	ug/L 98
22) 1,1-dichloroethene	6.02	96	147705	45.03	ug/L 94
23) acetone	6.05	58	13391	48.52	ug/L 92
24) allyl chloride	6.46	76	85345	35.54	ug/L # 59
25) acetonitrile	6.44	40	115220	436.10	ug/L # 58
26) iodomethane	6.27	142	299889	48.41	ug/L 99
27) carbon disulfide	6.38	76	455752	47.34	ug/L 99
28) methylene chloride	6.64	84	157523	45.33	ug/L 91
29) methyl acetate	6.44	74	23826	50.94	ug/L 92
30) 1-chloropropane	6.66	42	232053	39.13	ug/L 98
31) methyl tert butyl ether	6.89	73	405737	46.35	ug/L 98
32) trans-1,2-dichloroethene	6.95	96	140402	43.51	ug/L 94
33) di-isopropyl ether	7.41	45	490391	42.72	ug/L 97
34) 2-butanone	8.08	72	16704	45.43	ug/L # 71
35) 1,1-dichloroethane	7.45	63	258391	42.62	ug/L 98
36) chloroprene	7.55	53	208247	46.38	ug/L 97
37) acrylonitrile	6.92	53	269252	218.05	ug/L 100
38) vinyl acetate	7.43	86	22797	48.67	ug/L 99
39) ethyl tert-butyl ether	7.83	59	475652	46.64	ug/L 99
40) ethyl acetate	8.08	45	21027	42.13	ug/L 84
41) 2,2-dichloropropane	8.11	77	110438	43.08	ug/L 94
42) cis-1,2-dichloroethene	8.11	96	157278	44.87	ug/L 93

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54431.D
 Acq On : 12 Aug 2015 9:35 pm
 Operator : TOANP
 Sample : cc2289-50
 Misc : MS89499,V4B2297,w,,,1
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Aug 13 12:17:10 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.17	85	22590	47.46	ug/L	# 88
44) propionitrile	8.18	54	219489	440.69	ug/L	99
45) bromochloromethane	8.40	128	88131	48.99	ug/L	87
46) tetrahydrofuran	8.43	42	47797	39.94	ug/L	95
47) chloroform	8.44	85	168508	46.45	ug/L	95
48) T-BUTYL FORMATE	8.46	59	119052	51.34	ug/L	95
51) freon 113	5.96	151	98037	49.22	ug/L	100
52) methacrylonitrile	8.34	41	90451	41.12	ug/L	92
53) 1,1,1-trichloroethane	8.67	97	188639	49.18	ug/L	98
54) cyclohexane	8.74	84	177367	42.19	ug/L	95
55) iso-butyl alcohol	8.82	43	58940	430.59	ug/L	98
57) epichlorohydrin	10.63	57	76038	231.90	ug/L	99
58) n-butyl alcohol	9.53	56	221461	2381.82	ug/L	98
59) carbon tetrachloride	8.86	117	189343	49.54	ug/L	99
60) 1,1-dichloropropene	8.83	75	188316	44.35	ug/L	97
61) hexane	7.20	57	136418	42.30	ug/L	95
63) benzene	9.08	78	555115	44.41	ug/L	99
64) iso-octane	9.06	57	395449	39.39	ug/L	98
65) tert-amyl methyl ether	9.09	87	99518	50.08	ug/L	96
66) heptane	9.21	57	117599	52.96	ug/L	98
67) isopropyl acetate	8.99	61	61778	41.35	ug/L	96
68) 1,2-dichloroethane	9.10	62	195027	47.35	ug/L	95
69) trichloroethene	9.74	95	158154	47.01	ug/L	98
72) 2-nitropropane	10.50	41	61902	45.91	ug/L	# 84
73) 2-chloroethyl vinyl ether	10.50	63	466275	234.00	ug/L	97
74) methyl methacrylate	9.99	100	40103	50.03	ug/L	# 89
75) 1,2-dichloropropane	10.00	63	156488	43.58	ug/L	99
76) dibromomethane	10.16	93	102771	49.15	ug/L	96
77) methylcyclohexane	9.94	83	208042	48.09	ug/L	94
78) bromodichloromethane	10.28	83	217752	48.43	ug/L	97
79) cis-1,3-dichloropropene	10.72	75	263266	45.19	ug/L	98
81) 4-methyl-2-pentanone	10.82	58	63265	45.84	ug/L	97
82) toluene	11.09	92	368320	46.23	ug/L	100
83) 3-methyl-1-butanol	10.83	55	136861	888.19	ug/L	95
84) trans-1,3-dichloropropene	11.29	75	242636	47.56	ug/L	98
85) ethyl methacrylate	11.27	69	204459	48.44	ug/L	96
86) 1,1,2-trichloroethane	11.52	83	129964	46.86	ug/L	99
87) 2-hexanone	11.69	58	54986	43.63	ug/L	95
89) tetrachloroethene	11.69	164	181202	48.20	ug/L	99
90) 1,3-dichloropropane	11.71	76	236590	46.46	ug/L	97
91) butyl acetate	11.76	56	102266	44.02	ug/L	96
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	146511	429.81	ug/L	98
93) dibromochloromethane	11.99	129	194839	48.93	ug/L	100
94) 1,2-dibromoethane	12.15	107	165820	49.71	ug/L	100
96) chlorobenzene	12.63	112	438940	47.97	ug/L	98
97) 1,1,1,2-tetrachloroethane	12.70	131	168561	48.20	ug/L	99
98) ethylbenzene	12.69	91	720583	47.27	ug/L	99
99) m,p-xylene	12.80	106	551842	95.01	ug/L	98
100) o-xylene	13.25	106	289410	49.42	ug/L	97
101) styrene	13.27	104	481307	48.13	ug/L	97

Z
7.7.13

Quantitation Report (QT Reviewed)

Data Path : C:\MSDChem\1\DATA\
 Data File : 4B54431.D
 Acq On : 12 Aug 2015 9:35 pm
 Operator : TOANP
 Sample : cc2289-50
 Misc : MS89499,V4B2297,w,,,1
 ALS Vial : 27 Sample Multiplier: 1

Quant Time: Aug 13 12:17:10 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

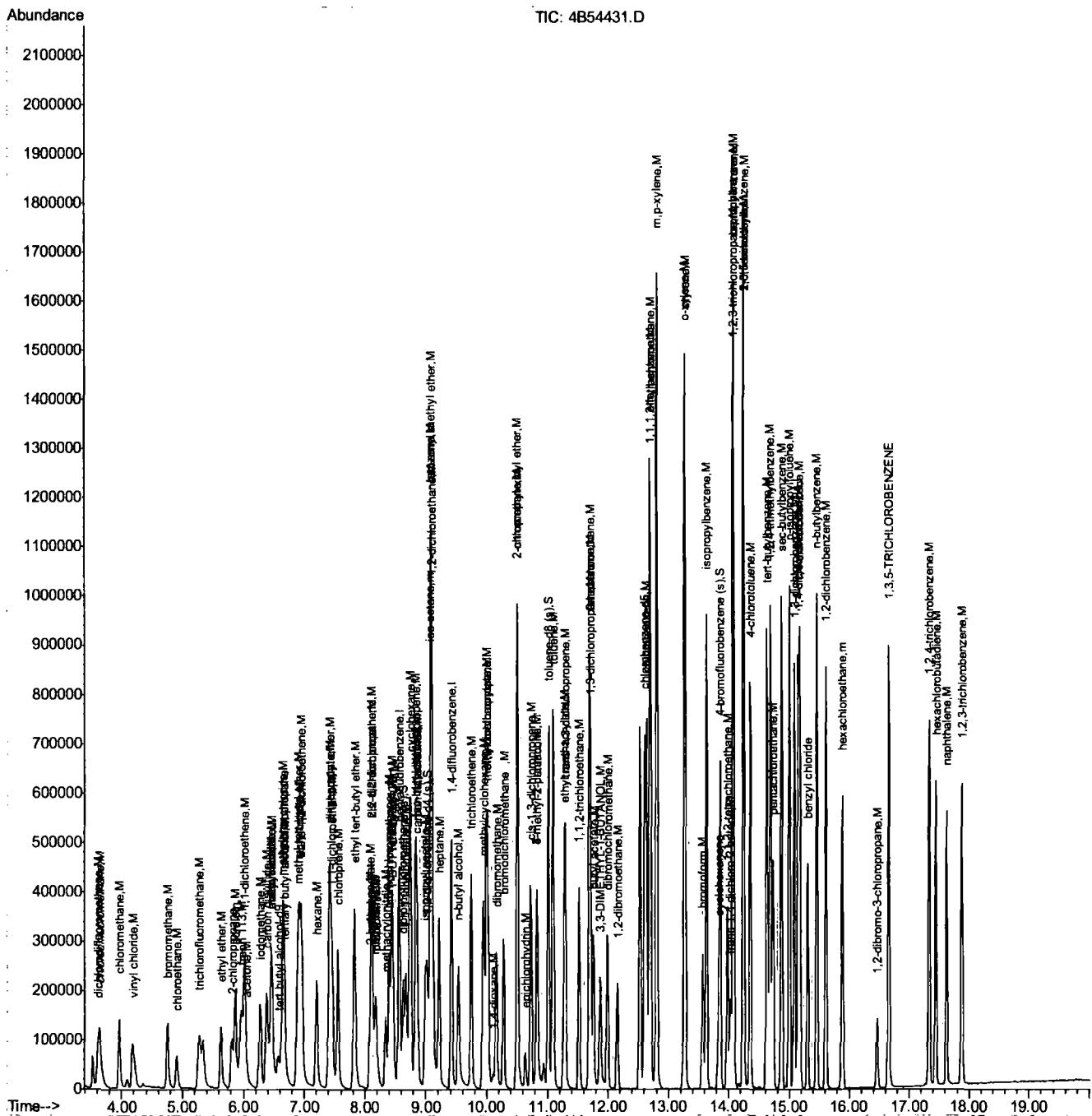
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
102) bromoform	13.57	173	151179	48.84	ug/L	99
104) isopropylbenzene	13.63	105	729179	45.67	ug/L	99
106) cyclohexanone	13.83	55	94188	206.75	ug/L	97
107) bromobenzene	14.07	156	224944	48.03	ug/L	94
108) 1,1,2,2-tetrachloroethane	13.98	83	218683	46.77	ug/L	99
109) trans-1,4-dichloro-2-buten	14.02	53	37297	31.02	ug/L	93
110) 1,2,3-trichloropropane	14.06	110	54081	48.81	ug/L	95
111) n-propylbenzene	14.08	91	824563	45.69	ug/L	99
113) 2-chlorotoluene	14.24	126	187490	45.84	ug/L	96
114) 4-chlorotoluene	14.35	91	531006	45.50	ug/L	99
115) 1,3,5-trimethylbenzene	14.24	105	622184	46.29	ug/L	100
116) tert-butylbenzene	14.62	119	560881	48.11	ug/L	99
117) pentachloroethane	14.72	167	118734	49.21	ug/L	98
118) 1,2,4-trimethylbenzene	14.68	105	616669	47.63	ug/L	99
119) sec-butylbenzene	14.87	105	827056	46.68	ug/L	100
120) 1,3-dichlorobenzene	15.08	146	419961	49.26	ug/L	100
121) p-isopropyltoluene	15.00	119	708109	47.81	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	410032	48.09	ug/L	99
123) benzyl chloride	15.30	91	367792	43.15	ug/L	99
124) 1,2-dichlorobenzene	15.61	146	415526	49.92	ug/L	98
126) n-butylbenzene	15.46	92	345918	47.09	ug/L	98
128) 1,2-dibromo-3-chloropropan	16.46	75	35456	47.25	ug/L	95
129) 1,3,5-TRICHLOROBENZENE	16.65	180	346698	52.22	ug/L	100
130) 1,2,4-trichlorobenzene	17.33	180	269455	50.67	ug/L	99
131) hexachlorobutadiene	17.44	225	167408	51.72	ug/L	98
132) naphthalene	17.62	128	470602	50.34	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	231318	51.57	ug/L	99
134) hexachloroethane	15.89	201	144789	47.15	ug/L	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54431.D
Acq On : 12 Aug 2015 9:35 pm
Operator : TOANP
Sample : cc2289-50
Misc : MS89499,V4B2297,w,,,1
ALS Vial : 27 Sample Multiplier: 1

Quant Time: Aug 13 12:17:10 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration





Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54455.D
Acq On : 13 Aug 2015 9:46 am
Operator : TOANP
Sample : cc2289-20
Misc : MS89470,V4B2298,w,,,1
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 13 16:12:44 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) tert butyl alcohol-d9	6.57	65	150330	500.00	ug/L	0.00
5) pentafluorobenzene	8.56	168	418083	50.00	ug/L	0.00
56) 1,4-difluorobenzene	9.41	114	472304	50.00	ug/L	0.00
88) chlorobenzene-d5	12.60	117	415435	50.00	ug/L	0.00
103) 1,4-dichlorobenzene-d4	15.14	152	226426	50.00	ug/L	0.00

System Monitoring Compounds

49) dibromofluoromethane (s)	8.62	113	148535	49.13	ug/L	0.00
Spiked Amount	50.000	Range	76 - 120	Recovery	=	98.26%
50) 1,2-dichloroethane-d4 (s)	9.01	65	157884	47.52	ug/L	0.00
Spiked Amount	50.000	Range	73 - 122	Recovery	=	95.04%
80) toluene-d8 (s)	11.01	98	546541	49.15	ug/L	0.00
Spiked Amount	50.000	Range	84 - 119	Recovery	=	98.30%
105) 4-bromofluorobenzene (s)	13.86	95	198566	50.23	ug/L	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	100.46%

Target Compounds

					Qvalue
2) tertiary butyl alcohol	6.67	59	39410	108.38	ug/L # 63
3) 1,4-dioxane	10.10	88	16946	504.06	ug/L 90
7) chlorodifluoromethane	3.64	51	76602	18.12	ug/L 94
8) dichlorodifluoromethane	3.60	85	94062	23.23	ug/L 97
10) chloromethane	3.97	52	30085	19.43	ug/L 88
11) vinyl chloride	4.17	62	79325	17.83	ug/L 96
12) bromomethane	4.77	94	57575	22.56	ug/L 96
13) chloroethane	4.92	64	38696	22.12	ug/L 98
15) trichlorofluoromethane	5.28	101	107238	25.51	ug/L 96
19) ethyl ether	5.63	74	29915	18.78	ug/L 96
20) 2-chloropropane	5.81	39	8541	16.85	ug/L 75
21) acrolein	5.86	56	98349	184.37	ug/L 94
22) 1,1-dichloroethene	6.03	96	67280	20.38	ug/L 95
23) acetone	6.05	58	6060	21.82	ug/L 94
24) allyl chloride	6.46	76	52998m	21.93	ug/L
25) acetonitrile	6.45	40	56108	211.03	ug/L # 69
26) iodomethane	6.28	142	118018	18.93	ug/L 99
27) carbon disulfide	6.38	76	181898	18.78	ug/L 98
28) methylene chloride	6.66	84	71545	20.46	ug/L 94
29) methyl acetate	6.44	74	9659	20.52	ug/L # 76
30) 1-chloropropane	6.65	42	107150	17.96	ug/L 95
31) methyl tert butyl ether	6.89	73	177561	20.16	ug/L 100
32) trans-1,2-dichloroethene	6.95	96	59900	18.45	ug/L 98
33) di-isopropyl ether	7.41	45	199338	17.26	ug/L 89
34) 2-butanone	8.08	72	7286	19.69	ug/L # 59
35) 1,1-dichloroethane	7.45	63	115504	18.93	ug/L 99
36) chloroprene	7.54	53	85663	18.96	ug/L 95
37) acrylonitrile	6.92	53	120628	97.08	ug/L 97
38) vinyl acetate	7.43	86	9093	19.29	ug/L 69
39) ethyl tert-butyl ether	7.82	59	193506	18.86	ug/L 99
40) ethyl acetate	8.08	45	9056	18.03	ug/L # 85
41) 2,2-dichloropropane	8.10	77	49026	19.01	ug/L 97
42) cis-1,2-dichloroethene	8.11	96	69104	19.59	ug/L 93

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54455.D
 Acq On : 13 Aug 2015 9:46 am
 Operator : TOANP
 Sample : cc2289-20
 Misc : MS89470,V4B2298,w,,,1
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 13 16:12:44 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) methylacrylate	8.16	85	9310	19.44	ug/L	97
44) propionitrile	8.18	54	98385	196.30	ug/L	87
45) bromochloromethane	8.40	128	38299	21.16	ug/L	86
46) tetrahydrofuran	8.42	42	21940	18.22	ug/L	96
47) chloroform	8.44	85	73808	20.22	ug/L	93
48) T-BUTYL FORMATE	8.46	59	45007	19.29	ug/L	98
51) freon 113	5.97	151	46751	23.33	ug/L	83
52) methacrylonitrile	8.34	41	39353	17.78	ug/L	95
53) 1,1,1-trichloroethane	8.67	97	76830	19.90	ug/L	97
54) cyclohexane	8.74	84	81645	19.30	ug/L	88
55) iso-butyl alcohol	8.82	43	30283	219.85	ug/L	97
57) epichlorohydrin	10.63	57	31372	96.75	ug/L	98
58) n-butyl alcohol	9.52	56	95710	1040.95	ug/L	97
59) carbon tetrachloride	8.85	117	79436	21.02	ug/L	96
60) 1,1-dichloropropene	8.83	75	85160	20.28	ug/L	99
61) hexane	7.20	57	73728	23.12	ug/L	97
63) benzene	9.08	78	244527	19.78	ug/L	98
64) iso-octane	9.06	57	192670	19.41	ug/L	99
65) tert-amyl methyl ether	9.09	87	39108	19.90	ug/L	90
66) heptane	9.21	57	65607	29.88	ug/L	89
67) isopropyl acetate	8.99	61	26711	18.08	ug/L	91
68) 1,2-dichloroethane	9.10	62	87600	21.51	ug/L	97
69) trichloroethene	9.73	95	68429	20.57	ug/L	98
72) 2-nitropropane	10.50	41	26812	20.11	ug/L	# 72
73) 2-chloroethyl vinyl ether	10.50	63	175981	89.31	ug/L	96
74) methyl methacrylate	9.98	100	17067	21.53	ug/L	# 84
75) 1,2-dichloropropane	10.00	63	67550	19.02	ug/L	99
76) dibromomethane	10.16	93	44549	21.55	ug/L	96
77) methylcyclohexane	9.94	83	98294	22.98	ug/L	95
78) bromodichloromethane	10.27	83	93624	21.06	ug/L	99
79) cis-1,3-dichloropropene	10.72	75	112370	19.51	ug/L	96
81) 4-methyl-2-pentanone	10.82	58	26880	19.69	ug/L	98
82) toluene	11.09	92	160732	20.40	ug/L	97
83) 3-methyl-1-butanol	10.83	55	62035	407.12	ug/L	92
84) trans-1,3-dichloropropene	11.29	75	101607	20.14	ug/L	98
85) ethyl methacrylate	11.27	69	83292	19.96	ug/L	96
86) 1,1,2-trichloroethane	11.52	83	55387	20.20	ug/L	96
87) 2-hexanone	11.69	58	23241	18.65	ug/L	87
89) tetrachloroethene	11.69	164	75007	20.78	ug/L	99
90) 1,3-dichloropropane	11.71	76	101948	20.85	ug/L	99
91) butyl acetate	11.76	56	38982	17.48	ug/L	96
92) 3,3-DIMETHYL-1-BUTANOL	11.87	57	65631	200.54	ug/L	97
93) dibromochloromethane	11.99	129	82270	21.52	ug/L	94
94) 1,2-dibromoethane	12.15	107	69485	21.70	ug/L	98
96) chlorobenzene	12.63	112	185858	21.15	ug/L	97
97) 1,1,1,2-tetrachloroethane	12.69	131	72080	21.47	ug/L	99
98) ethylbenzene	12.68	91	304923	20.84	ug/L	98
99) m,p-xylene	12.80	106	237472	42.58	ug/L	100
100) o-xylene	13.25	106	121567	21.62	ug/L	99
101) styrene	13.27	104	198537	20.68	ug/L	96

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
 Data File : 4B54455.D
 Acq On : 13 Aug 2015 9:46 am
 Operator : TOANP
 Sample : cc2289-20
 Misc : MS89470,V4B2298,w,,,1
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 13 16:12:44 2015
 Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
 Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 QLast Update : Mon Aug 10 08:19:32 2015
 Response via : Initial Calibration

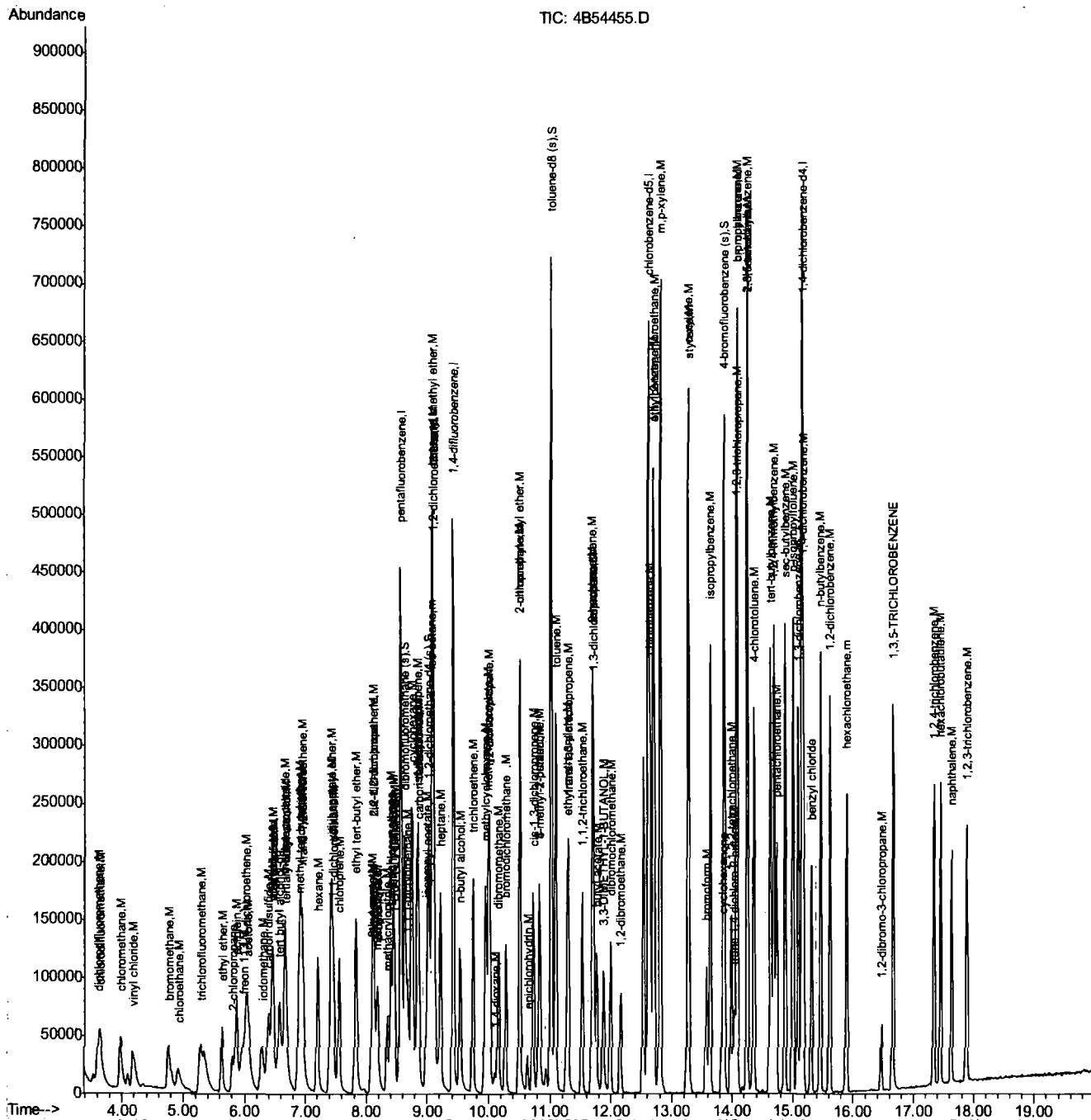
Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
102) bromoform	13.57	173	62053	20.88	ug/L	98
104) isopropylbenzene	13.63	105	309817	21.55	ug/L	98
106) cyclohexanone	13.82	55	39660	96.70	ug/L	97
107) bromobenzene	14.07	156	94748	22.47	ug/L	94
108) 1,1,2,2-tetrachloroethane	13.98	83	91641	21.77	ug/L	99
109) trans-1,4-dichloro-2-butene	14.02	53	16407	15.16	ug/L	96
110) 1,2,3-trichloropropane	14.05	110	23675	23.74	ug/L	89
111) n-propylbenzene	14.08	91	338064	20.81	ug/L	98
113) 2-chlorotoluene	14.24	126	78090	21.21	ug/L	96
114) 4-chlorotoluene	14.35	91	217681	20.72	ug/L	98
115) 1,3,5-trimethylbenzene	14.24	105	257573	21.29	ug/L	98
116) tert-butylbenzene	14.62	119	230775	21.99	ug/L	98
117) pentachloroethane	14.72	167	54920	25.28	ug/L	97
118) 1,2,4-trimethylbenzene	14.68	105	252059	21.63	ug/L	99
119) sec-butylbenzene	14.87	105	334274	20.96	ug/L	99
120) 1,3-dichlorobenzene	15.08	146	169306	22.06	ug/L	99
121) p-isopropyltoluene	15.00	119	283683	21.28	ug/L	99
122) 1,4-dichlorobenzene	15.17	146	161965	21.10	ug/L	99
123) benzyl chloride	15.30	91	157573	20.54	ug/L	98
124) 1,2-dichlorobenzene	15.61	146	166428	22.21	ug/L	98
126) n-butylbenzene	15.46	92	129900	19.64	ug/L	99
128) 1,2-dibromo-3-chloropropan	16.46	75	14402	21.32	ug/L	93
129) 1,3,5-TRICHLOROBENZENE	16.65	180	130956	21.91	ug/L	100
130) 1,2,4-trichlorobenzene	17.33	180	95782	20.01	ug/L	99
131) hexachlorobutadiene	17.44	225	69583	23.88	ug/L	100
132) naphthalene	17.62	128	171611	20.39	ug/L	99
133) 1,2,3-trichlorobenzene	17.88	180	83561	20.70	ug/L	99
134) hexachloroethane	15.89	201	61838	22.37	ug/L	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Quantitation Report (QT Reviewed)

Data Path : C:\MSDCHEM\1\DATA\
Data File : 4B54455.D
Acq On : 13 Aug 2015 9:46 am
Operator : TOANP
Sample : cc2289-20
Misc : MS89470,V4B2298,w,,,1
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Aug 13 16:12:44 2015
Quant Method : C:\MSDCHEM\1\METHODS\M4B2289.M
Quant Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
QLast Update : Mon Aug 10 08:19:32 2015
Response via : Initial Calibration



Manual Integration Approval Summary

Page 1 of 1

Sample Number: V4B2298-CC2289 **Method:** SW846 8260C
Lab FileID: 4B54455.D **Analyst approved:** 08/13/15 16:31 Dong, Mei
Injection Time: 08/13/15 09:46 **Supervisor approved:** 08/13/15 16:53 Kanya Veerawat

Parameter	CAS	Sig#	R.T. (min.)	Reason
Allyl chloride	107-05-1		6.46	Poor instrument integration

7.7.14.1

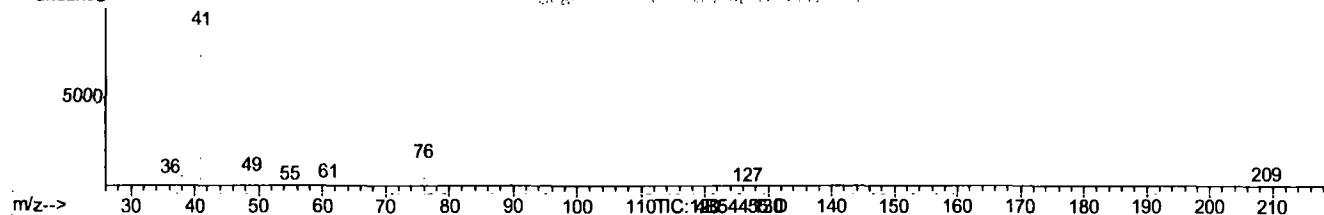
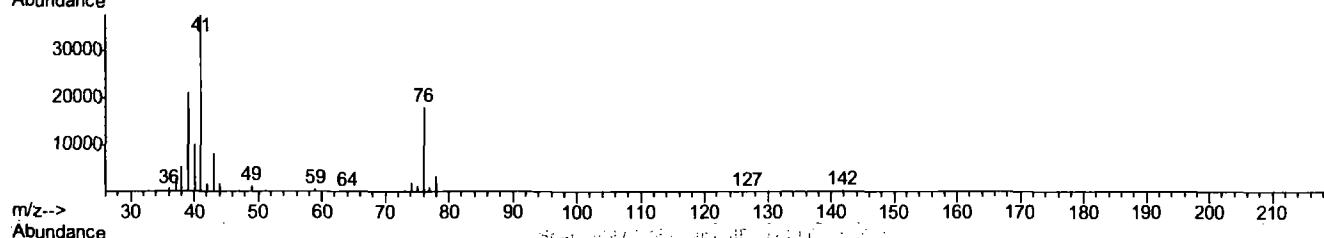
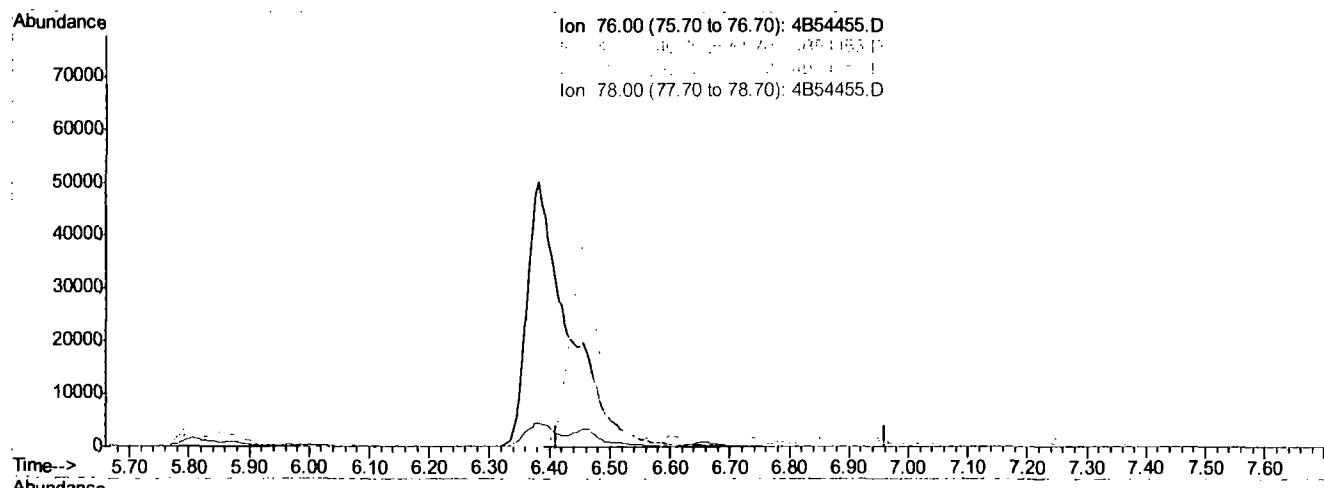


Quantitation Report (Qedit)

Data File : C:\MSDCHEM\1\DATA\4B54455.D Vial: 2
 Acq On : 13 Aug 2015 9:46 am Operator: TOANP
 Sample : cc2289-20 Inst : MS4B
 Misc : MS89470,V4B2298,w,,,,,1 Multiplr: 1.00
 MS Integration Params: RTEINT.P

Quant Time: Aug 13 10:27:09 2015 Results File: M4B2289.RES

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 Last Update : Mon Aug 10 08:19:32 2015
 Response via : Multiple Level Calibration



(24) allyl chloride (M)

6.46min 0.00ug/L

response 0

Ion	Exp%	Act%
76.00	100	0.00
41.00	208.30	0.00#
39.00	110.00	0.00#
78.00	18.00	0.00

Quantitation Report (Qedit)

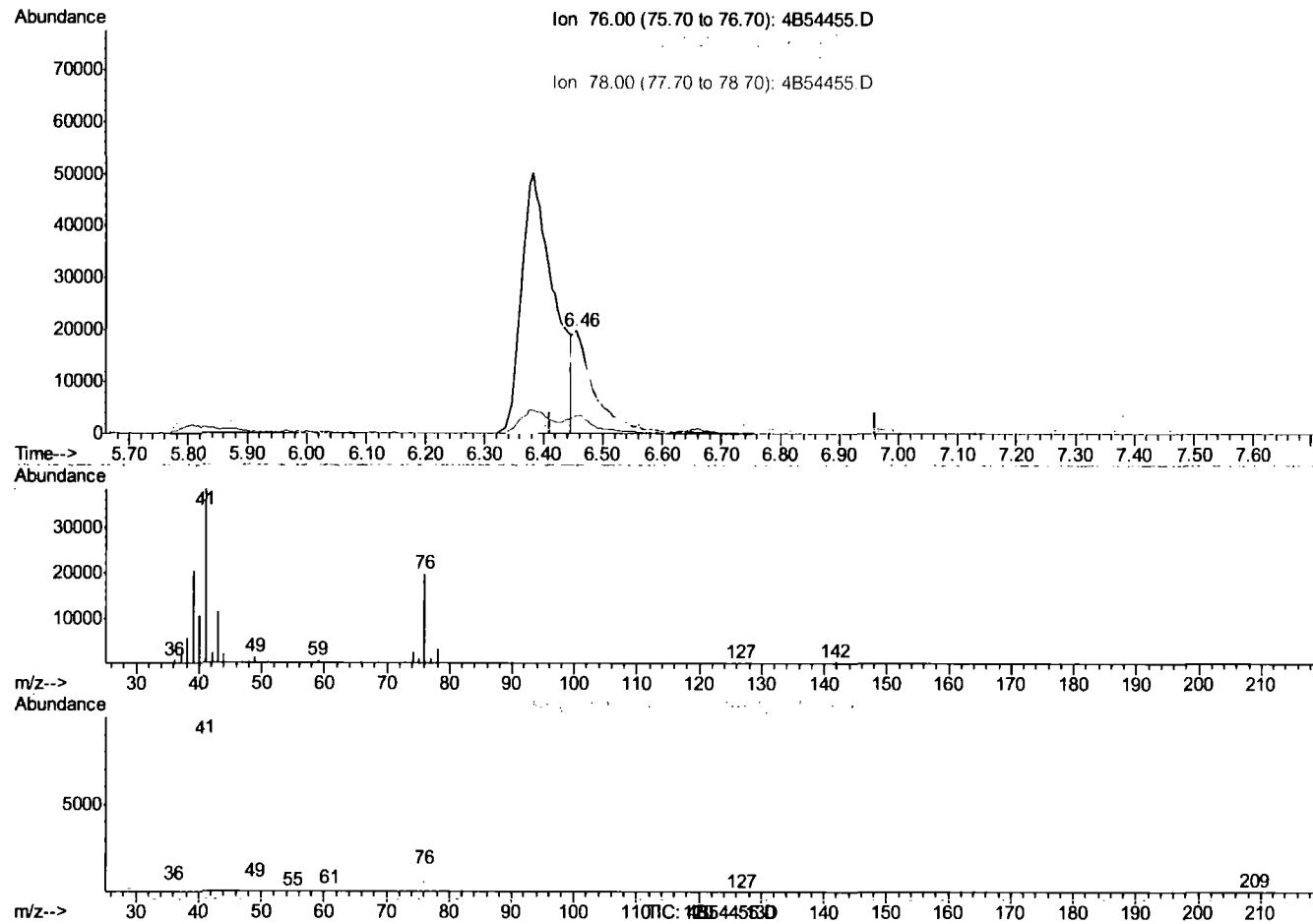
Data File : C:\MSDCHEM\1\DATA\4B54455.D
 Acq On : 13 Aug 2015 9:46 am
 Sample : cc2289-20
 Misc : MS89470,V4B2298,w,,,1
 MS Integration Params: RTEINT.P

Vial: 2
 Operator: TOANP
 Inst : MS4B
 Multiplr: 1.00

Quant Time: Aug 13 10:27:09 2015

Results File: M4B2289.RES

Method : C:\MSDCHEM\1\METHODS\M4B2289.M (RTE Integrator)
 Title : Method SW846 V8260C, ZB624 60mx0.25mmx1.4um
 Last Update : Mon Aug 10 08:19:32 2015
 Response via : Multiple Level Calibration



(24) allyl chloride (M)

6.46min 21.93ug/L m

response 52998

Ion	Exp%	Act%
76.00	100	100
41.00	208.30	194.73
39.00	110.00	103.89
78.00	18.00	17.39



VOLATILE ANALYSIS LOG

Date: 8/15/15

Standard Data

Lot #	Description	Cone.
W015-7114-93.4	A	100 ppm
W015-7114-72.45	B	100 ppm
W015-7114-96.16	C	100 ppm
W015-7114-52	E. Acrylate	100 ppm
W015-7114-34.46	HEx	100 ppm

Standard Data

Lot #	Description	Cone.
W015-7114-95.4	Ext A	100 ppm
W015-7114-73.12	Ext B	100 ppm
W015-7114-94.5	Ext C	100 ppm
W015-7114-71.4	Hex	100 ppm
W015-7114- Ext E. Acrylate		100 ppm

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature:

Date: 8/15/15

R	Data File	Sample ID	Test	M Vial T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt. (uL)	Secondary dilution	L +	I S	S U	Status (Data)	Comments	pH* <2
	4B54241	APF1												OK	
	54242	IC2289-0-2	Y610	W			5							OK 100/500ml	
	54243	IC2289-0-5	V	W			5							OK 100/100ml	
	54244	IC2289-1	V	W			5							OK 100/100ml	
	54245	IC2289-2	V	W			5							OK 100/100ml	
	54246	IC2289-5	V	W			5							OK 100/100ml	
	54247	IC2289-10	V	W			5							OK 100/100ml	
	54248	IC2289-20	V	W			5							OK 100/100ml	
	54249	IC2289-50	V	W			5							OK 100/100ml	
	54250	IC2289-100	V	W			5							OK 100/100ml	
	54251	IC2289-200	V	W			5							OK 100/100ml	
	54252	IB													
	54253	SB													
	54254	IC2289-50	8210	W			5							Ext A 100/100ml	
	54255	IB													

MTX = Matrix Designate W for water, S for soil, O for oil. L+ =Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt= Volume (ML) or Weight (g); MOH amt.= volume (uL) extract injected * IF pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9

Rev. Date: 2/14/2007

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ACCUTEST.

VOLATILE ANALYSIS LOG

Date: 8/12/15

Batch ID: V4B 2296

Print Analyst Name: DCIN Pham

Analyst Signature:

Columns: 25m x 0.25mm x 1.4µm

Method 840C

Initial Cal. Method V4B 2296

Standard Data		
Lot #	Description	Conc.
V015-104-13-25	A	100 ppm
V015-104-12-13	B	100 ppm
V015-104-106-2	C	100 ppm
V015-104-107-9	AC	1000 ppm
216814	pH paper exp/06/15/17	100 ppm

Standard Data		
Lot #	Description	Conc.
V015-104-14-17	Ext A	100 ppm
V015-104-14-18	Ext B	100 ppm
V015-104-14-1	Ext C	100 ppm
V015-104-107-1	Ext AC	1000 ppm
V015-104-109-9	ILS	100 ppm

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature:

Date: 8/12/15

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt. (uL)	Secondary dilution	L + I S U	S U	Status (Data)	Comments	pH* <2
	4B54405	NPB											OK	
	54406	LC D284-10					5						OK	
	54407	LC1049-10					5						1	
	54408	IS					5						OK	
	54409	MB					5						OK	
	54410	BS					5						OK	
	54411	IA					5						OK	
K	54412	TC100-2	893745 TC100	G	1		5		10x				OK	
K	54413	TC960-3	893930 TC960	G	2		5		10x				OK	
K	54414	TC960-3	✓ 89470	G	2		5/10		10x				OK	
F	54415	TC100-1	893745 TC100	G	2		5/10		10x				OK	
	54416	TC1107-3	89470	G	3		5		10x				OK	
	54417	TC1004-1	89494 TC1004	G	4		5		10x				OK	
	54418	TC1004-2	89494 TC1004	G	4		5		10x				OK	
	54419	TC1004-3	89470 TC1004	G	4		5		10x				OK	
	54420	TC1107-3 mg	89470 SL	G									OK	
	54421	TC1107-3 mg	89470 SL	G									OK	

MTX = Matrix Designate W for water, S for soil, O for oil. L+ =Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt.= volume (uL) extract injected * IF pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

1

Form: OR001-9

Rev. Date: 2/14/2007

7.8.2



ACCUTEST.

VOLATILE ANALYSIS LOG

Date: 8/14/05

Standard Data

Standard Data

Lot #	Description	Conc.
101C		

Batch ID: V46 2296

Print Analyst Name: Tawn Phoenix

Analyst Signature:

Columns: 23624 160m x 25mm x 1.4 fm

Method

Initial Cal. Method M 48 2259

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EO44.

Supervisor Signature:

[Signature] Date: 2/19/15

MTX = Matrix Designate W for water, S for soil, O for oil. **L+** = Library Search. **IS** = Internal Standard Area. **SU** = Surrogate.

Sample Amt = Volume (mL) or Weight (g): MOH amt = volume (mL) extract injected * If pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error.

Form: OR001-9

Rev. Date: 2/14/2007

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VOLATILE ANALYSIS LOG

Batch ID: V4B2297

Date: 8/18/15

Standard Data

Lot #	Description	Cone.
V01T-WM - 03.15	A	100 ppm
V01T-WM - 72.53	B	100 ppm
V01T-WM - 106.1	C	100 ppm
V01T-WM - 101.9	AC	100 ppm
21664 pH pepti exp 6/15/17		

Standard Data

Lot #	Description	Cone.
V01T-WM - 95.17	EXT A	100 ppm
V01T-WM - 071311	EXT B	100 ppm
V01T-WM - 111.1	EXT C	100 ppm
V01T-WM - 107.1	EXT AV	1000 ppm
V01T-WM - 89	1000 ppm	1000 ppm

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature:

Date: 8/18/15

R	Data File	Sample ID	Test	M/Vial T X	ALS #	Samp. Amt (ml or g)	MOH amt. (ul)	Secondary dilution	L + S U	I	S	U	Status (Data)	Comments	pH <2
	54434-20	APB				"								Ac	
	54431	TC2289-4				5								✓	A,B,C, ACV
	54432	DB												✓	✓
	54433	MB				5								✓	✓
	54434	BS				6								✓	✓
	54435	TC869-1 in TC2289-4	84344	G	2	5/10		10x						reviewed ext A, B, C, ACV	✓
	54436	TC869-1mL	✓	G	2	5/10		10x						✓	✓
	54437	TB												✓	
	54438	TC 869-5 TC2289-4	84344	G	2	5		1x						✓	
	54439	TC869-6	✓	G	2	5		1x						✓	
	54440	TC 869-1	✓	G	2	5/10		10x						✓	RR1x +com
	54441	TC 869-3	✓	G	1	5		1x						✓	
	54442	TC 869-4	✓	G	2	5		1x						✓	
	54443	TC 869-2	✓	G	2	5		1x						✓	
	54444	TC 869-2	✓	G	2	5/100		10x						+4334443	✓
	54445	TC1024-1 TC1024-1	84444	E	1	5		1x						844444 TC1024-1 other not	✓
	54446	TC1024-4 TC1024-2	✓	G	1	5		1x						✓	

MTX = Matrix Designate W for water, S for soil, O for oil. L+ =Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt= volume (ul) extract injected * If pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9

Rev. Date: 2/14/2007

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7.8.3



ACCUTEST.

VOLATILE ANALYSIS LOG

Date: 8/12/15

Standard Data

Lot #	Description	Conc.

Standard Data

Lot #	Description	Conc.
5		
part		
0		

Batch ID: V4B2297

Print Analyst Name: Toucan Draw

Analyst Signature: Jes

Columns: 2x6x14 (60mm x 25mm x 1.4cm)

Method

Initial Cal. Method M4B 2289

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EOA044.

Supervisor Signature:

Date: 8/18/16

MTX = Matrix Designate W for water, S for soil, O for oil. **L+** = Library Search. **IS = Internal Standard Area.** **SU = Surrogate.**

Sample Designations: W for water, S for soil, G for oil. L = Laboratory Search. LS = Laboratory Standard Area. LR = Laboratory Reference.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error.

Form: OR001-9

Rev. Date: 2/14/2007



VOLATILE ANALYSIS LOG

Date: 8/13/05

Standard Data

Lot #	Description	Conc.
VDT 2M1	45-17	60ppm
VDT 2M1	120-1	100 ppm
VDT 2M1	111-1	100 ppm
VDT 2M1	107-1	1000 ppm
VDT 2M1	92-2	100 ppm

Standard Data

Lot #	Description	Conc.
VDT 2M1	93-25	A
VDT 2M1	74-53	B
VDT 2M1	106-2	C
VDT 2M1	101-1	A/B
VDT 2M1	89	D/E

Batch ID: V4B2297/V2B2298

Print Analyst Name: Jean Jobe Williams

Analyst Signature:

Columns: ZB621

Method: V82602

Initial Cal. Method 4B2281

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature:

Date: 8/14/2005

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt (uL)	Secondary dilution	L + S	I U	S U	Status (Data)	Comments	pH <2
	4B54454	BFB												9.00ml	
	54455	CC 2209-20												10ml 50ml NIST on AIB C1100	
	54456	IB													
	54457	MB2													
	54458	JC B60-3 PUP	TLL20+100	G	2										
	54459	JC B60-4 MS	✓	G	1									21.0ml 143ml AIB C1100	
	54460	IB													
	54461	MB1													
	54462	BS													
DIL	54463	JC979-B	B9406 TLL20	G W	4		(100uL)/50		500X					25ml 50ml 500 E4 PUP, Hex 100 NOT different from 100ml after centrifugation	
	54464	IC1045-1S	B9444 SL9	G W	3	5			1X						
	54465	JC1024-i	B9499 TLL20+100	G W	5	5			1X	+ W					
	54466	JC1024-2	✓	G W	5	5			1X	+ W					
	54467	JC1045-15MS	B9444 SL9	G W	1	5			1X						
	54468	JC1045-15MSD	✓	G W	4	5			1X						
DIL	54469	JC 979-B	B9406 TLL20	G W	5		(200uL)/50		250X					NOT USC Vial difference	
	54470	JC1024-3	B9444 TLL20+100	G W	5	5			1X						

MTX = Matrix Designate W for water, S for soil, O for oil. L+ = Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt = volume (uL) extract injected * IF pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9

Rev. Date: 2/14/2007

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VOLATILE ANALYSIS LOG

Date: 8/13/01

Standard Data

Lot #	Description	Conc.
VDR 2M	95-19	6x18
VDR 2M	120-1	6x18
VDR 2M	111-1	6x18
VDR 2M	107-1	6x18
VDR 2M	92-2	6x18

Standard Data

Lot #	Description	Conc.
VDR 2M	93-25	A
VDR 2M	71-53	B
VDR 2M	106-2	C
VDR 2M	101-1	F18
VDR 2M	89	I18

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature:

Date: 8/14/2001

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ml or g)	MOH amt. (ul)	Secondary dilution	L + S	I	S U	Status (Date)	Comments	pH* <2
	4854454	BFB												4.00ml	
	54455	JL 2289-20												10ml 50ml 11B, C, D, E	
	54456	IB													
	54457	MB2													
	54458	JL 1260-3 PUP	TLL20+TBL	G	2										
	54459	JL 1261-4 ms		G	1									21.6ml 145ml A,B,C,D,E	
	54460	IB													
	54461	MB1													
	54462	BS													
DIL R	54463	JL 979-B	89406 TLL20	G W	4		(100ul)/50		500X					25ml 5ml F14 Alight, Max 18ml NOT USE Dilution difference seen in alight tll20	
	54464	JL 1045-1 S	89449 SL9	G W	3	5			1X						
R	54465	JL 1024-i	89499 TLL20+TBL	G W	5	5			1X	+ W					
R	54466	JL 1024-2		G W	5	5			1X	+ W					
	54467	JL 1045-15 ms	89449 SL9	G W	1	5			1X						
	54468	JL 1045-15MSD		G W	1	5			1X						
DIL R	54469	JL 979-S	89406 TLL20	G W	3		(200ul) ₅₀		250X					NOT USE O Vial difference	
	54470	JL 1024-3	89499 TBL	G W	5	5			1X						

MTX = Matrix Designate W for water, S for soil, O for oil. L+ =Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt.= volume (ul) extract injected * IF pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9

Rev. Date: 2/14/2007

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VOLATILE ANALYSIS LOG

Batch ID: V2B2298

Date: 8/13/05

Print Analyst Name: ZB624 H am?

Standard Data

Lot #	Description	Conc.

Standard Data

Lot #	Description	Conc.
	ref pg 13	

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature:

Date: 8/13/2005

7.84

Z

R	Data File	Sample ID	Test	M T X	Vial #	ALS #	Samp. Amt (ML or g)	MOH amt. (uL)	Secondary dilution	L + S	I S U	Status (Data)	Comments	pH < 2
			894499	G	4		5		1x	✓		OK		✓
54471	JC1045-1S	T6261792	894499	G	1		5		1x	✓		OK		✓
54472	JC1045-1	SL9	894499	G	1		5		1x	✓		OK		✓
54473	JC1045-2	V	894499	G	1		5		1x	✓		OK		✓
54474	JC1045-3	V	894499	G	1		5		1x	✓		OK		✓
54475	JC1045-4	V	894499	G	2		5		1x	✓		OK		✓
54476	JC1045-5	V	894499	G	2		5		1x	✓		OK	8:15 pm	✓
54477	BFB2											OK	8:43 pm	
54478	CC2298-50											✓	OK	25ml 50ml AIBCAW
54479	1B											✓	clip	
54480	mB2											✓	OK	
54481	JC1045-6	894499	894499	G	1		5		1x	✓		OK		✓
54482	JC1045-7	V	894499	G	2		5		1x	✓		OK		✓
54483	JC1045-8	V	894499	G	2		5		1x	✓		OK		✓
54484	JC1045-9	V	894499	G	2		5		1x	✓		OK		✓
54485	JC1045-10	V	894499	G	1		5		1x	✓		OK		✓
54486	JC1045-11	V	894499	G	1		5		1x	✓		OK		✓
54487	JC1045-12	V	894499	G	2		5		1x	✓		OK		✓

MTX = Matrix Designate W for water, S for soil, O for oil. L+ =Library Search. IS = Internal Standard Area. SU = Surrogate.

Sample Amt = Volume (ML) or Weight (g); MOH amt.= volume (uL) extract injected * IF pH > 2, comment on sample result.

All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analyst's correction error

Form: OR001-9

Rev. Date: 2/14/2007

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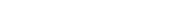
VOLATILE ANALYSIS LOG

Date: 8/13/15

Standard Data

Standard Data

Manually integrated chromatographic peaks in the following reportable files have been reviewed and verified to comply with the criteria of Accutest SOP EQA044.

Supervisor Signature: 

Date: 8/14/2015

Initial Cal. Method MgA₂ 2289

Columns: ZB62

Method V82401

MTX = Matrix Designate W for water, S for soil, O for oil. **L+** =Library Search. **IS = Internal Standard Area.** **SU = Surrogate.**

Sample Amt = Volume (ML) or Weight (g). MOH amt = volume (uL) extract injected * 1E pH > 2 comment on sample result

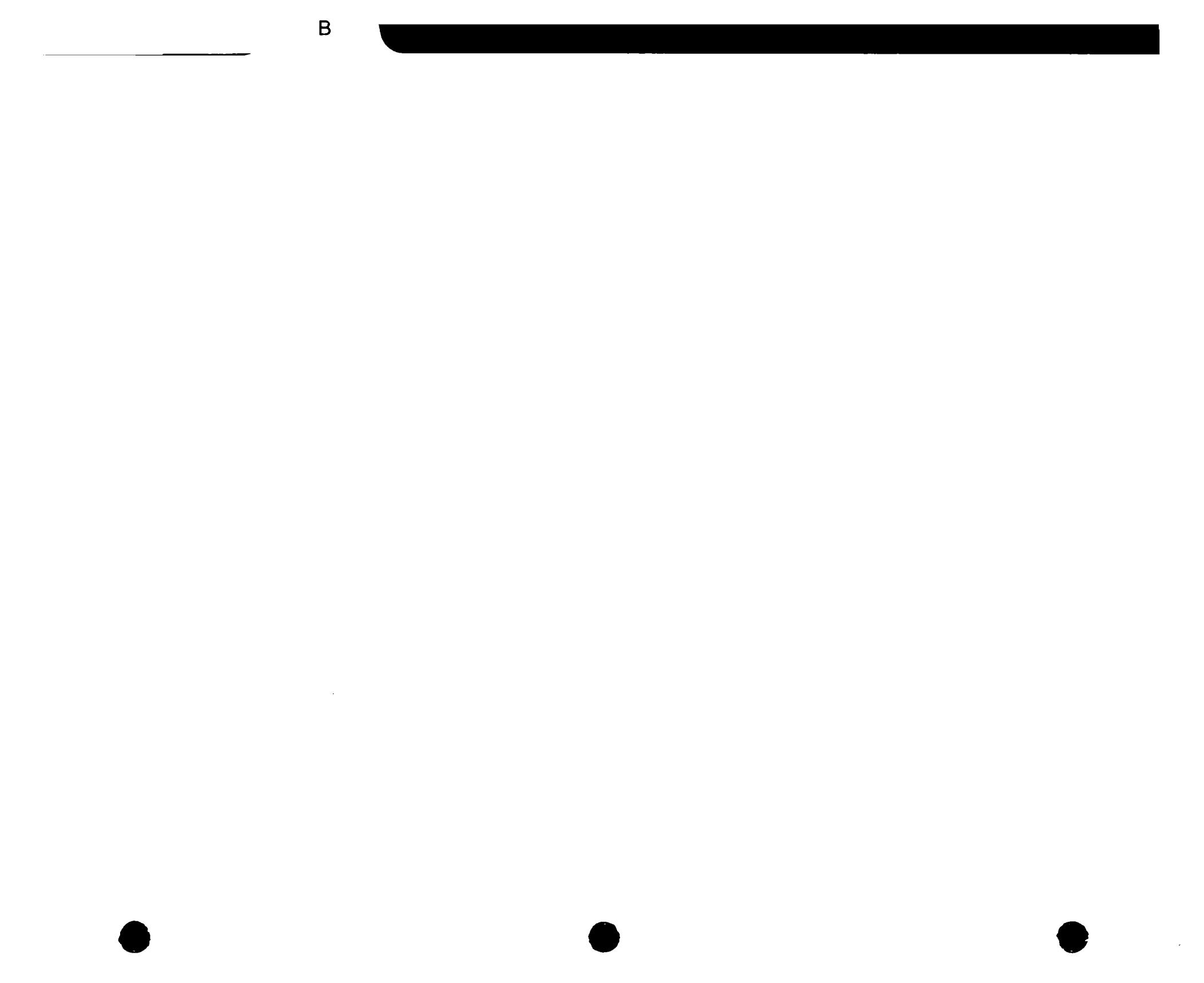
All strike outs must be initialed, dated and reason code applied as follows: 1 = reviewer correction error; 2 = transcription error; 3 = computer miscalculation; 4 = analysis correction error.

Form: OB001 a

Form. OR001-9
Rev. Date: 3/14/2007

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B



B

Appendix B

Third Quarter 2015 Effluent Air Laboratory Analytical Reports

7/30/2015

Mr. Peter Hollatz
AECOM Environment
27755 Diehl Road
Suite 100
Warrenville IL 60555

Project Name: UTC-SER-9/10

Project #: 60339110
Workorder #: 1507260

Dear Mr. Peter Hollatz

The following report includes the data for the above referenced project for sample(s) received on 7/17/2015 at Air Toxics Ltd.

The data and associated QC analyzed by Modified TO-14A are compliant with the project requirements or laboratory criteria with the exception of the deviations noted in the attached case narrative.

Thank you for choosing Air Toxics Ltd. for your air analysis needs. Air Toxics Ltd. is committed to providing accurate data of the highest quality. Please feel free to contact the Project Manager: Ausha Scott at 916-985-1000 if you have any questions regarding the data in this report.

Regards,



Ausha Scott

Project Manager

A Eurofins Lancaster Laboratories Company

Eurofins Air Toxics, Inc.

180 Blue Ravine Road, Suite B
Folsom, CA 95630

T | 916-985-1000
F | 916-985-1020
www.airtoxics.com



Air Toxics

WORK ORDER #: 1507260

Work Order Summary

CLIENT: Mr. Peter Hollatz
 AECOM Environment
 27755 Diehl Road
 Suite 100
 Warrenville, IL 60555

BILL TO: Accounts Payable-Warrenville
 AECOM Environment
 27755 Diehl Road
 Suite 100
 Warrenville, IL 60555

PHONE: 630-836-1700

P.O. #: 60339110

FAX: 630-836-1711

PROJECT #: 60339110 UTC-SER-9/10

DATE RECEIVED: 07/17/2015

CONTACT: Ausha Scott

DATE COMPLETED: 07/30/2015

<u>FRACTION #</u>	<u>NAME</u>	<u>TEST</u>	<u>RECEIPT</u>	<u>VAC./PRES.</u>	<u>FINAL PRESSURE</u>
01A	P1SVE-EFFC2-0716/15	Modified TO-14A	3.9 "Hg	14.7 psi	
02A	P1SVE-EFFC3-0716/15	Modified TO-14A	5.5 "Hg	15 psi	
03A	P1SVE-EFFC1-0716/15	Modified TO-14A	3.3 "Hg	14.9 psi	
04A	P2SVE-EFFC4-0716/15	Modified TO-14A	5.1 "Hg	15.1 psi	
05A	P2SVE-EFFC5-0716/15	Modified TO-14A	4.1 "Hg	14.9 psi	
06A	Lab Blank	Modified TO-14A	NA	NA	
07A	CCV	Modified TO-14A	NA	NA	
08A	LCS	Modified TO-14A	NA	NA	
08AA	LCSD	Modified TO-14A	NA	NA	

CERTIFIED BY:

DATE: 07/30/15

Technical Director

Certification numbers: AZ Licensure AZ0775, NJ NELAP - CA016, NY NELAP - 11291,

TX NELAP - T104704343-14-7, UT NELAP CA009332014-5, VA NELAP - 460197, WA NELAP - C935

Name of Accreditation Body: NELAP/ORELAP (Oregon Environmental Laboratory Accreditation Program)

Accreditation number: CA300005, Effective date: 10/18/2014, Expiration date: 10/17/2015.

Eurofins Air Toxics Inc.. certifies that the test results contained in this report meet all requirements of the NELAC standards

This report shall not be reproduced, except in full, without the written approval of Eurofins Air Toxics, Inc.

180 BLUE RAVINE ROAD, SUITE B FOLSOM, CA - 95630

(916) 985-1000 . (800) 985-5955 . FAX (916) 985-1020

LABORATORY NARRATIVE

Modified TO-14A
AECOM Environment
Workorder# 1507260

Five 1 Liter Summa Canister samples were received on July 17, 2015. The laboratory performed analysis via modified EPA Method TO-14A using GC/MS in the full scan mode.

This workorder was independently validated prior to submittal using 'USEPA National Functional Guidelines' as generally applied to the analysis of volatile organic compounds in air. A rules-based, logic driven, independent validation engine was employed to assess completeness, evaluate pass/fail of relevant project quality control requirements and verification of all quantified amounts.

Method modifications taken to run these samples are summarized in the table below. Specific project requirements may over-ride the ATL modifications. Please note that TO-14A was validated for specially treated canisters, and the use of Tedlar bags for sample collection is outside the scope of the method.

Requirement	TO-14A	ATL Modifications
Initial Calibration criteria	RSD</=30%	Follow TO-15 requirements of RSD</=30% with two compounds allowed out to </=40%RSD.
BFB absolute abundance criteria	Within 10% of that from previous day	CCV internal standard area counts are compared to ICAL, corrective action when recovery is less than 60%.
Blank acceptance criteria	<0.20 ppbv	<Reporting Limit
Sample Drying System	Nafion Dryer	Multibed hydrophobic sorbent
BFB ion abundance criteria	Ion abundance listed in Table 4 of TO-14A	Follow ion abundance criteria listed in Method TO-15

Receiving Notes

There were no receiving discrepancies.

Analytical Notes

There were no analytical discrepancies.

Definition of Data Qualifying Flags

Nine qualifiers may have been used on the data analysis sheets and indicates as follows:

B - Compound present in laboratory blank greater than reporting limit (background subtraction not performed).

J - Estimated value.

E - Exceeds instrument calibration range.

S - Saturated peak.

Q - Exceeds quality control limits.

U - Compound analyzed for but not detected above the reporting limit, LOD, or MDL value. See data page for project specific U-flag definition.

UJ- Non-detected compound associated with low bias in the CCV

N - The identification is based on presumptive evidence.

C - Estimated calculation due to estimated sampling rate.

File extensions may have been used on the data analysis sheets and indicates

as follows:

a-File was requantified

b-File was quantified by a second column and detector

r1-File was requantified for the purpose of reissue



Air Toxics

Summary of Detected Compounds
MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Client Sample ID: P1SVE-EFFC2-0716/15

Lab ID#: 1507260-01A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.2	1.7	4.6	6.9
1,1-Dichloroethane	1.2	17	4.6	71
cis-1,2-Dichloroethene	1.2	4.2	4.6	17
1,1,1-Trichloroethane	1.2	64	6.3	350
Trichloroethene	1.2	18	6.2	98
Tetrachloroethene	1.2	8.6	7.8	58

Client Sample ID: P1SVE-EFFC3-0716/15

Lab ID#: 1507260-02A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.2	2.3	4.9	9.0
1,1-Dichloroethane	1.2	6.5	5.0	26
cis-1,2-Dichloroethene	1.2	19	4.9	77
1,1,1-Trichloroethane	1.2	180	6.7	980
Trichloroethene	1.2	12	6.6	63
Tetrachloroethene	1.2	6.0	8.4	41

Client Sample ID: P1SVE-EFFC1-0716/15

Lab ID#: 1507260-03A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.1	3.1	4.5	12
1,1-Dichloroethane	1.1	32	4.6	130
cis-1,2-Dichloroethene	1.1	6.7	4.5	27
1,1,1-Trichloroethane	1.1	110	6.2	590
Trichloroethene	1.1	19	6.1	100
Tetrachloroethene	1.1	170	7.7	1200

Client Sample ID: P2SVE-EFFC4-0716/15

Lab ID#: 1507260-04A



Summary of Detected Compounds MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Client Sample ID: P2SVE-EFFC4-0716/15

Lab ID#: 1507260-04A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.2	4.0	4.8	16
1,1-Dichloroethane	1.2	7.7	4.9	31
cis-1,2-Dichloroethene	1.2	13	4.8	52
1,1,1-Trichloroethane	1.2	270	6.6	1500
Trichloroethene	1.2	27	6.6	140
Tetrachloroethene	1.2	130	8.3	880

Client Sample ID: P2SVE-EFFC5-0716/15

Lab ID#: 1507260-05A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.2	3.5	4.6	14
1,1-Dichloroethane	1.2	22	4.7	89
cis-1,2-Dichloroethene	1.2	12	4.6	47
1,1,1-Trichloroethane	1.2	260	6.4	1400
Trichloroethene	1.2	14	6.3	78
Tetrachloroethene	1.2	54	7.9	370



Air Toxics

Client Sample ID: P1SVE-EFFC2-0716/15

Lab ID#: 1507260-01A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072008	Date of Collection: 7/16/15 9:00:00 AM		
Dil. Factor:	2.30	Date of Analysis: 7/20/15 03:40 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.2	Not Detected	2.9	Not Detected
Chloroethane	4.6	Not Detected	12	Not Detected
1,1-Dichloroethene	1.2	1.7	4.6	6.9
Acetone	12	Not Detected	27	Not Detected
Methylene Chloride	12	Not Detected	40	Not Detected
trans-1,2-Dichloroethene	1.2	Not Detected	4.6	Not Detected
1,1-Dichloroethane	1.2	17	4.6	71
2-Butanone (Methyl Ethyl Ketone)	4.6	Not Detected	14	Not Detected
cis-1,2-Dichloroethene	1.2	4.2	4.6	17
Chloroform	1.2	Not Detected	5.6	Not Detected
1,1,1-Trichloroethane	1.2	64	6.3	350
Carbon Tetrachloride	1.2	Not Detected	7.2	Not Detected
Benzene	1.2	Not Detected	3.7	Not Detected
1,2-Dichloroethane	1.2	Not Detected	4.6	Not Detected
Trichloroethene	1.2	18	6.2	98
Toluene	1.2	Not Detected	4.3	Not Detected
1,1,2-Trichloroethane	1.2	Not Detected	6.3	Not Detected
Tetrachloroethene	1.2	8.6	7.8	58
Ethyl Benzene	1.2	Not Detected	5.0	Not Detected
m,p-Xylene	1.2	Not Detected	5.0	Not Detected
o-Xylene	1.2	Not Detected	5.0	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	116	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	82	70-130



Air Toxics

Client Sample ID: P1SVE-EFFC3-0716/15

Lab ID#: 1507260-02A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072009	Date of Collection:	7/16/15 9:25:00 AM
Dil. Factor:	2.47	Date of Analysis:	7/20/15 04:17 PM
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)
Vinyl Chloride	1.2	Not Detected	3.2
Chloroethane	4.9	Not Detected	13
1,1-Dichloroethene	1.2	2.3	4.9
Acetone	12	Not Detected	29
Methylene Chloride	12	Not Detected	43
trans-1,2-Dichloroethene	1.2	Not Detected	4.9
1,1-Dichloroethane	1.2	6.5	5.0
2-Butanone (Methyl Ethyl Ketone)	4.9	Not Detected	14
cis-1,2-Dichloroethene	1.2	19	4.9
Chloroform	1.2	Not Detected	6.0
1,1,1-Trichloroethane	1.2	180	6.7
Carbon Tetrachloride	1.2	Not Detected	7.8
Benzene	1.2	Not Detected	3.9
1,2-Dichloroethane	1.2	Not Detected	5.0
Trichloroethene	1.2	12	6.6
Toluene	1.2	Not Detected	4.6
1,1,2-Trichloroethane	1.2	Not Detected	6.7
Tetrachloroethene	1.2	6.0	8.4
Ethyl Benzene	1.2	Not Detected	5.4
m,p-Xylene	1.2	Not Detected	5.4
o-Xylene	1.2	Not Detected	5.4

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	113	70-130
Toluene-d8	104	70-130
4-Bromofluorobenzene	82	70-130



Air Toxics

Client Sample ID: P1SVE-EFFC1-0716/15

Lab ID#: 1507260-03A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072010	Date of Collection: 7/16/15 9:55:00 AM		
Dil. Factor:	2.26	Date of Analysis: 7/20/15 04:40 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.1	Not Detected	2.9	Not Detected
Chloroethane	4.5	Not Detected	12	Not Detected
1,1-Dichloroethene	1.1	3.1	4.5	12
Acetone	11	Not Detected	27	Not Detected
Methylene Chloride	11	Not Detected	39	Not Detected
trans-1,2-Dichloroethene	1.1	Not Detected	4.5	Not Detected
1,1-Dichloroethane	1.1	32	4.6	130
2-Butanone (Methyl Ethyl Ketone)	4.5	Not Detected	13	Not Detected
cis-1,2-Dichloroethene	1.1	6.7	4.5	27
Chloroform	1.1	Not Detected	5.5	Not Detected
1,1,1-Trichloroethane	1.1	110	6.2	590
Carbon Tetrachloride	1.1	Not Detected	7.1	Not Detected
Benzene	1.1	Not Detected	3.6	Not Detected
1,2-Dichloroethane	1.1	Not Detected	4.6	Not Detected
Trichloroethene	1.1	19	6.1	100
Toluene	1.1	Not Detected	4.2	Not Detected
1,1,2-Trichloroethane	1.1	Not Detected	6.2	Not Detected
Tetrachloroethene	1.1	170	7.7	1200
Ethyl Benzene	1.1	Not Detected	4.9	Not Detected
m,p-Xylene	1.1	Not Detected	4.9	Not Detected
o-Xylene	1.1	Not Detected	4.9	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	112	70-130
Toluene-d8	104	70-130
4-Bromofluorobenzene	82	70-130



Air Toxics

Client Sample ID: P2SVE-EFFC4-0716/15

Lab ID#: 1507260-04A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072013	Date of Collection: 7/16/15 10:05:00 AM		
Dil. Factor:	2.44	Date of Analysis: 7/20/15 06:17 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.2	Not Detected	3.1	Not Detected
Chloroethane	4.9	Not Detected	13	Not Detected
1,1-Dichloroethene	1.2	4.0	4.8	16
Acetone	12	Not Detected	29	Not Detected
Methylene Chloride	12	Not Detected	42	Not Detected
trans-1,2-Dichloroethene	1.2	Not Detected	4.8	Not Detected
1,1-Dichloroethane	1.2	7.7	4.9	31
2-Butanone (Methyl Ethyl Ketone)	4.9	Not Detected	14	Not Detected
cis-1,2-Dichloroethene	1.2	13	4.8	52
Chloroform	1.2	Not Detected	6.0	Not Detected
1,1,1-Trichloroethane	1.2	270	6.6	1500
Carbon Tetrachloride	1.2	Not Detected	7.7	Not Detected
Benzene	1.2	Not Detected	3.9	Not Detected
1,2-Dichloroethane	1.2	Not Detected	4.9	Not Detected
Trichloroethene	1.2	27	6.6	140
Toluene	1.2	Not Detected	4.6	Not Detected
1,1,2-Trichloroethane	1.2	Not Detected	6.6	Not Detected
Tetrachloroethene	1.2	130	8.3	880
Ethyl Benzene	1.2	Not Detected	5.3	Not Detected
m,p-Xylene	1.2	Not Detected	5.3	Not Detected
o-Xylene	1.2	Not Detected	5.3	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	114	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	82	70-130



Air Toxics

Client Sample ID: P2SVE-EFFC5-0716/15

Lab ID#: 1507260-05A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072014	Date of Collection: 7/16/15 10:10:00 AM		
Dil. Factor:	2.33	Date of Analysis: 7/20/15 06:41 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.2	Not Detected	3.0	Not Detected
Chloroethane	4.7	Not Detected	12	Not Detected
1,1-Dichloroethene	1.2	3.5	4.6	14
Acetone	12	Not Detected	28	Not Detected
Methylene Chloride	12	Not Detected	40	Not Detected
trans-1,2-Dichloroethene	1.2	Not Detected	4.6	Not Detected
1,1-Dichloroethane	1.2	22	4.7	89
2-Butanone (Methyl Ethyl Ketone)	4.7	Not Detected	14	Not Detected
cis-1,2-Dichloroethene	1.2	12	4.6	47
Chloroform	1.2	Not Detected	5.7	Not Detected
1,1,1-Trichloroethane	1.2	260	6.4	1400
Carbon Tetrachloride	1.2	Not Detected	7.3	Not Detected
Benzene	1.2	Not Detected	3.7	Not Detected
1,2-Dichloroethane	1.2	Not Detected	4.7	Not Detected
Trichloroethene	1.2	14	6.3	78
Toluene	1.2	Not Detected	4.4	Not Detected
1,1,2-Trichloroethane	1.2	Not Detected	6.4	Not Detected
Tetrachloroethene	1.2	54	7.9	370
Ethyl Benzene	1.2	Not Detected	5.0	Not Detected
m,p-Xylene	1.2	Not Detected	5.0	Not Detected
o-Xylene	1.2	Not Detected	5.0	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	114	70-130
Toluene-d8	101	70-130
4-Bromofluorobenzene	86	70-130



Air Toxics

Client Sample ID: Lab Blank

Lab ID#: 1507260-06A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072005	Date of Collection:	NA	
Dil. Factor:	1.00	Date of Analysis:	7/20/15 10:39 AM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	0.50	Not Detected	1.3	Not Detected
Chloroethane	2.0	Not Detected	5.3	Not Detected
1,1-Dichloroethene	0.50	Not Detected	2.0	Not Detected
Acetone	5.0	Not Detected	12	Not Detected
Methylene Chloride	5.0	Not Detected	17	Not Detected
trans-1,2-Dichloroethene	0.50	Not Detected	2.0	Not Detected
1,1-Dichloroethane	0.50	Not Detected	2.0	Not Detected
2-Butanone (Methyl Ethyl Ketone)	2.0	Not Detected	5.9	Not Detected
cis-1,2-Dichloroethene	0.50	Not Detected	2.0	Not Detected
Chloroform	0.50	Not Detected	2.4	Not Detected
1,1,1-Trichloroethane	0.50	Not Detected	2.7	Not Detected
Carbon Tetrachloride	0.50	Not Detected	3.1	Not Detected
Benzene	0.50	Not Detected	1.6	Not Detected
1,2-Dichloroethane	0.50	Not Detected	2.0	Not Detected
Trichloroethene	0.50	Not Detected	2.7	Not Detected
Toluene	0.50	Not Detected	1.9	Not Detected
1,1,2-Trichloroethane	0.50	Not Detected	2.7	Not Detected
Tetrachloroethene	0.50	Not Detected	3.4	Not Detected
Ethyl Benzene	0.50	Not Detected	2.2	Not Detected
m,p-Xylene	0.50	Not Detected	2.2	Not Detected
o-Xylene	0.50	Not Detected	2.2	Not Detected

Container Type: NA - Not Applicable

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	114	70-130
Toluene-d8	104	70-130
4-Bromofluorobenzene	84	70-130



Air Toxics

Client Sample ID: CCV

Lab ID#: 1507260-07A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072002	Date of Collection: NA
Dil. Factor:	1.00	Date of Analysis: 7/20/15 08:40 AM

Compound	%Recovery
Vinyl Chloride	119
Chloroethane	120
1,1-Dichloroethene	101
Acetone	123
Methylene Chloride	125
trans-1,2-Dichloroethene	104
1,1-Dichloroethane	116
2-Butanone (Methyl Ethyl Ketone)	111
cis-1,2-Dichloroethene	98
Chloroform	106
1,1,1-Trichloroethane	99
Carbon Tetrachloride	96
Benzene	112
1,2-Dichloroethane	112
Trichloroethene	99
Toluene	106
1,1,2-Trichloroethane	110
Tetrachloroethene	93
Ethyl Benzene	105
m,p-Xylene	105
o-Xylene	102

Container Type: NA - Not Applicable

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	117	70-130
Toluene-d8	108	70-130
4-Bromofluorobenzene	87	70-130



Air Toxics

Client Sample ID: LCS

Lab ID#: 1507260-08A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072003	Date of Collection: NA
Dil. Factor:	1.00	Date of Analysis: 7/20/15 09:14 AM
Compound	%Recovery	Method Limits
Vinyl Chloride	124	70-130
Chloroethane	127	70-130
1,1-Dichloroethene	103	70-130
Acetone	124	70-130
Methylene Chloride	125	70-130
trans-1,2-Dichloroethene	90	70-130
1,1-Dichloroethane	116	70-130
2-Butanone (Methyl Ethyl Ketone)	113	70-130
cis-1,2-Dichloroethene	107	70-130
Chloroform	108	70-130
1,1,1-Trichloroethane	98	70-130
Carbon Tetrachloride	98	70-130
Benzene	114	70-130
1,2-Dichloroethane	114	70-130
Trichloroethene	101	70-130
Toluene	108	70-130
1,1,2-Trichloroethane	112	70-130
Tetrachloroethene	94	70-130
Ethyl Benzene	109	70-130
m,p-Xylene	106	70-130
o-Xylene	106	70-130

Container Type: NA - Not Applicable

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	116	70-130
Toluene-d8	109	70-130
4-Bromofluorobenzene	89	70-130



Air Toxics

Client Sample ID: LCSD

Lab ID#: 1507260-08AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17072004	Date of Collection: NA
Dil. Factor:	1.00	Date of Analysis: 7/20/15 09:49 AM
Compound	%Recovery	Method Limits
Vinyl Chloride	125	70-130
Chloroethane	127	70-130
1,1-Dichloroethene	104	70-130
Acetone	124	70-130
Methylene Chloride	128	70-130
trans-1,2-Dichloroethene	92	70-130
1,1-Dichloroethane	115	70-130
2-Butanone (Methyl Ethyl Ketone)	114	70-130
cis-1,2-Dichloroethene	109	70-130
Chloroform	109	70-130
1,1,1-Trichloroethane	98	70-130
Carbon Tetrachloride	98	70-130
Benzene	112	70-130
1,2-Dichloroethane	112	70-130
Trichloroethene	101	70-130
Toluene	106	70-130
1,1,2-Trichloroethane	108	70-130
Tetrachloroethene	92	70-130
Ethyl Benzene	104	70-130
m,p-Xylene	103	70-130
o-Xylene	104	70-130

Container Type: NA - Not Applicable

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	105	70-130
Toluene-d8	109	70-130
4-Bromofluorobenzene	88	70-130

10/6/2015

Mr. Peter Hollatz
AECOM Environment
27755 Diehl Road
Suite 100
Warrenville IL 60555

Project Name: UTC-HS-PLT 1

Project #: 60339110
Workorder #: 1509395

Dear Mr. Peter Hollatz

The following report includes the data for the above referenced project for sample(s) received on 9/23/2015 at Air Toxics Ltd.

The data and associated QC analyzed by Modified TO-14A are compliant with the project requirements or laboratory criteria with the exception of the deviations noted in the attached case narrative.

Thank you for choosing Eurofins Air Toxics Inc. for your air analysis needs. Eurofins Air Toxics Inc. is committed to providing accurate data of the highest quality. Please feel free the Project Manager: Ausha Scott at 916-985-1000 if you have any questions regarding the data in this report.

Regards,



Ausha Scott

Project Manager

A Eurofins Lancaster Laboratories Company

Eurofins Air Toxics, Inc.

180 Blue Ravine Road, Suite B
Folsom, CA 95630

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www.airtoxics.com



Air Toxics

WORK ORDER #: 1509395

Work Order Summary

CLIENT:	Mr. Peter Hollatz AECOM Environment 27755 Diehl Road Suite 100 Warrenville, IL 60555	BILL TO:	Accounts Payable-Warrenville AECOM Environment 27755 Diehl Road Suite 100 Warrenville, IL 60555
PHONE:	630-836-1700	P.O. #	58743ACM
FAX:	630-836-1711	PROJECT #	60339110 UTC-HS-PLT 1
DATE RECEIVED:	09/23/2015	CONTACT:	Ausha Scott
DATE COMPLETED:	10/05/2015		

<u>FRACTION #</u>	<u>NAME</u>	<u>TEST</u>	<u>RECEIPT VAC./PRES.</u>	<u>FINAL PRESSURE</u>
01A	P1SVE-EFFC1-092215	Modified TO-14A	8 "Hg	14.8 psi
01AA	P1SVE-EFFC1-092215 Lab Duplicate	Modified TO-14A	8 "Hg	14.8 psi
02A	P2SVE-EFFC4-092215	Modified TO-14A	2.2 "Hg	14.9 psi
03A	P2SVE-EFFC5-092215	Modified TO-14A	5.1 "Hg	15.3 psi
04A	P1SVE-EFFC2-092215	Modified TO-14A	3.5 "Hg	14.6 psi
05A	P1SVE-EFFC2-092215 DUP	Modified TO-14A	4.9 "Hg	14.8 psi
06A	P1SVE-EFFC3-092215	Modified TO-14A	3.9 "Hg	14.9 psi
07A	FIELD BLANK-092215	Modified TO-14A	2.2 "Hg	15 psi
08A	Lab Blank	Modified TO-14A	NA	NA
09A	CCV	Modified TO-14A	NA	NA
10A	LCS	Modified TO-14A	NA	NA
10AA	LCSD	Modified TO-14A	NA	NA

CERTIFIED BY:

DATE: 10/06/15

Technical Director

Certification numbers: AZ Licensure AZ0775, NJ NELAP - CA016, NY NELAP - 11291,
 TX NELAP - T104704343-14-7, UT NELAP CA009332014-5, VA NELAP - 460197, WA NELAP - C935
 Name of Accreditation Body: NELAP/ORELAP (Oregon Environmental Laboratory Accreditation Program)

Accreditation number: CA300005, Effective date: 10/18/2014, Expiration date: 10/17/2015.

Eurofins Air Toxics Inc.. certifies that the test results contained in this report meet all requirements of the NELAC standards

This report shall not be reproduced, except in full, without the written approval of Eurofins Air Toxics, Inc.

180 BLUE RAVINE ROAD, SUITE B FOLSOM, CA - 9563
 (916) 985-1000 . (800) 985-5955 . FAX (916) 985-1020

**LABORATORY NARRATIVE
Modified TO-14A
AECOM Environment
Workorder# 1509395**

Seven 1 Liter Summa Canister samples were received on September 23, 2015. The laboratory performed analysis via modified EPA Method TO-14A using GC/MS in the full scan mode.

This workorder was independently validated prior to submittal using 'USEPA National Functional Guidelines' as generally applied to the analysis of volatile organic compounds in air. A rules-based, logic driven, independent validation engine was employed to assess completeness, evaluate pass/fail of relevant project quality control requirements and verification of all quantified amounts.

Method modifications taken to run these samples are summarized in the table below. Specific project requirements may over-ride the ATL modifications. Please note that TO-14A was validated for specially treated canisters, and the use of Tedlar bags for sample collection is outside the scope of the method.

Requirement	TO-14A	ATL Modifications
Initial Calibration criteria	RSD</=30%	Follow TO-15 requirements of RSD</=30% with two compounds allowed out to </=40%RSD.
BFB absolute abundance criteria	Within 10% of that from previous day	CCV internal standard area counts are compared to ICAL, corrective action when recovery is less than 60%.
Blank acceptance criteria	<0.20 ppbv	<Reporting Limit
Sample Drying System	Nafion Dryer	Multibed hydrophobic sorbent
BFB ion abundance criteria	Ion abundance listed in Table 4 of TO-14A	Follow ion abundance criteria listed in Method TO-15

Receiving Notes

The Chain of Custody (COC) information for sample P1SVE-EFFC1-092215 did not match the information on the canister with regard to canister identification. The client was notified of the discrepancy and the information on the canister was used to process and report the sample.

Analytical Notes

Dilution was performed on sample P1SVE-EFFC3-092215 due to the presence of high level target species.

The field blank sample FIELD BLANK-092215 has reportable levels of target compounds present.

Definition of Data Qualifying Flags

Nine qualifiers may have been used on the data analysis sheets and indicates as follows:

B - Compound present in laboratory blank greater than reporting limit (background subtraction not performed).

J - Estimated value.

E - Exceeds instrument calibration range.

S - Saturated peak.

Q - Exceeds quality control limits.

U - Compound analyzed for but not detected above the reporting limit, LOD, or MDL value. See data page for project specific U-flag definition.

UJ- Non-detected compound associated with low bias in the CCV

N - The identification is based on presumptive evidence.

C - Estimated calculation due to estimated sampling rate.

File extensions may have been used on the data analysis sheets and indicates as follows:

a-File was requantified

b-File was quantified by a second column and detector

r1-File was requantified for the purpose of reissue



Summary of Detected Compounds MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Client Sample ID: P1SVE-EFFC1-092215

Lab ID#: 1509395-01A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethane	1.4	29	5.5	120
cis-1,2-Dichloroethene	1.4	5.6	5.4	22
1,1,1-Trichloroethane	1.4	150	7.5	840
Trichloroethene	1.4	20	7.4	110
Tetrachloroethene	1.4	250	9.3	1700
Ethyl Benzene	1.4	2.2	5.9	9.5
m,p-Xylene	1.4	3.4	5.9	15

Client Sample ID: P1SVE-EFFC1-092215 Lab Duplicate

Lab ID#: 1509395-01AA

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethane	1.4	28	5.5	110
cis-1,2-Dichloroethene	1.4	5.2	5.4	20
1,1,1-Trichloroethane	1.4	150	7.5	820
Trichloroethene	1.4	21	7.4	110
Tetrachloroethene	1.4	250	9.3	1700
Ethyl Benzene	1.4	2.0	5.9	8.7
m,p-Xylene	1.4	3.6	5.9	16

Client Sample ID: P2SVE-EFFC4-092215

Lab ID#: 1509395-02A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.1	2.1	4.3	8.2
1,1-Dichloroethane	1.1	6.3	4.4	26
cis-1,2-Dichloroethene	1.1	11	4.3	45
1,1,1-Trichloroethane	1.1	200	5.9	1100
Trichloroethene	1.1	36	5.8	190
Tetrachloroethene	1.1	200	7.4	1400



Summary of Detected Compounds

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Client Sample ID: P2SVE-EFFC5-092215

Lab ID#: 1509395-03A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.2	1.2	4.9	5.0
1,1-Dichloroethane	1.2	4.2	5.0	17
cis-1,2-Dichloroethene	1.2	2.4	4.9	9.7
1,1,1-Trichloroethane	1.2	150	6.7	810
Trichloroethene	1.2	11	6.6	61
Tetrachloroethene	1.2	47	8.3	320

Client Sample ID: P1SVE-EFFC2-092215

Lab ID#: 1509395-04A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.1	3.4	4.5	14
1,1-Dichloroethane	1.1	210	4.6	860
cis-1,2-Dichloroethene	1.1	9.6	4.5	38
1,1,1-Trichloroethane	1.1	450	6.2	2500
Trichloroethene	1.1	30	6.1	160
Tetrachloroethene	1.1	28	7.7	190

Client Sample ID: P1SVE-EFFC2-092215 DUP

Lab ID#: 1509395-05A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	1.2	3.8	4.8	15
1,1-Dichloroethane	1.2	210	4.8	840
cis-1,2-Dichloroethene	1.2	9.2	4.8	37
1,1,1-Trichloroethane	1.2	440	6.5	2400
Trichloroethene	1.2	31	6.4	170
Tetrachloroethene	1.2	19	8.1	130

Client Sample ID: P1SVE-EFFC3-092215

Lab ID#: 1509395-06A



Summary of Detected Compounds
MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

Client Sample ID: P1SVE-EFFC3-092215

Lab ID#: 1509395-06A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1-Dichloroethene	2.3	2.6	9.2	10
1,1-Dichloroethane	2.3	11	9.4	43
cis-1,2-Dichloroethene	2.3	10	9.2	41
1,1,1-Trichloroethane	2.3	530	13	2900
Trichloroethene	2.3	14	12	73
Tetrachloroethene	2.3	18	16	120

Client Sample ID: FIELD BLANK-092215

Lab ID#: 1509395-07A

Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
1,1,1-Trichloroethane	1.1	1.2	5.9	6.7



Air Toxics

Client Sample ID: P1SVE-EFFC1-092215

Lab ID#: 1509395-01A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092809	Date of Collection: 9/22/15 9:05:00 AM		
Dil. Factor:	2.74	Date of Analysis: 9/28/15 02:49 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.4	Not Detected	3.5	Not Detected
Chloroethane	5.5	Not Detected	14	Not Detected
1,1-Dichloroethene	1.4	Not Detected	5.4	Not Detected
Acetone	14	Not Detected	32	Not Detected
Methylene Chloride	14	Not Detected	48	Not Detected
trans-1,2-Dichloroethene	1.4	Not Detected	5.4	Not Detected
1,1-Dichloroethane	1.4	29	5.5	120
2-Butanone (Methyl Ethyl Ketone)	5.5	Not Detected	16	Not Detected
cis-1,2-Dichloroethene	1.4	5.6	5.4	22
Chloroform	1.4	Not Detected	6.7	Not Detected
1,1,1-Trichloroethane	1.4	150	7.5	840
Carbon Tetrachloride	1.4	Not Detected	8.6	Not Detected
Benzene	1.4	Not Detected	4.4	Not Detected
1,2-Dichloroethane	1.4	Not Detected	5.5	Not Detected
Trichloroethene	1.4	20	7.4	110
Toluene	1.4	Not Detected	5.2	Not Detected
1,1,2-Trichloroethane	1.4	Not Detected	7.5	Not Detected
Tetrachloroethene	1.4	250	9.3	1700
Ethyl Benzene	1.4	2.2	5.9	9.5
m,p-Xylene	1.4	3.4	5.9	15
o-Xylene	1.4	Not Detected	5.9	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	99	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	93	70-130



Air Toxics

Client Sample ID: P1SVE-EFFC1-092215 Lab Duplicate

Lab ID#: 1509395-01AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092810	Date of Collection:	9/22/15 9:05:00 AM
Dil. Factor:	2.74	Date of Analysis:	9/28/15 03:12 PM
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)
Vinyl Chloride	1.4	Not Detected	3.5
Chloroethane	5.5	Not Detected	14
1,1-Dichloroethene	1.4	Not Detected	5.4
Acetone	14	Not Detected	32
Methylene Chloride	14	Not Detected	48
trans-1,2-Dichloroethene	1.4	Not Detected	5.4
1,1-Dichloroethane	1.4	28	5.5
2-Butanone (Methyl Ethyl Ketone)	5.5	Not Detected	16
cis-1,2-Dichloroethene	1.4	5.2	5.4
Chloroform	1.4	Not Detected	6.7
1,1,1-Trichloroethane	1.4	150	7.5
Carbon Tetrachloride	1.4	Not Detected	8.6
Benzene	1.4	Not Detected	4.4
1,2-Dichloroethane	1.4	Not Detected	5.5
Trichloroethene	1.4	21	7.4
Toluene	1.4	Not Detected	5.2
1,1,2-Trichloroethane	1.4	Not Detected	7.5
Tetrachloroethene	1.4	250	9.3
Ethyl Benzene	1.4	2.0	5.9
m,p-Xylene	1.4	3.6	5.9
o-Xylene	1.4	Not Detected	5.9

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	102	70-130
Toluene-d8	103	70-130
4-Bromofluorobenzene	96	70-130



Air Toxics

Client Sample ID: P2SVE-EFFC4-092215

Lab ID#: 1509395-02A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092811	Date of Collection: 9/22/15 9:25:00 AM		
Dil. Factor:	2.17	Date of Analysis: 9/28/15 03:41 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.1	Not Detected	2.8	Not Detected
Chloroethane	4.3	Not Detected	11	Not Detected
1,1-Dichloroethene	1.1	2.1	4.3	8.2
Acetone	11	Not Detected	26	Not Detected
Methylene Chloride	11	Not Detected	38	Not Detected
trans-1,2-Dichloroethene	1.1	Not Detected	4.3	Not Detected
1,1-Dichloroethane	1.1	6.3	4.4	26
2-Butanone (Methyl Ethyl Ketone)	4.3	Not Detected	13	Not Detected
cis-1,2-Dichloroethene	1.1	11	4.3	45
Chloroform	1.1	Not Detected	5.3	Not Detected
1,1,1-Trichloroethane	1.1	200	5.9	1100
Carbon Tetrachloride	1.1	Not Detected	6.8	Not Detected
Benzene	1.1	Not Detected	3.5	Not Detected
1,2-Dichloroethane	1.1	Not Detected	4.4	Not Detected
Trichloroethene	1.1	36	5.8	190
Toluene	1.1	Not Detected	4.1	Not Detected
1,1,2-Trichloroethane	1.1	Not Detected	5.9	Not Detected
Tetrachloroethene	1.1	200	7.4	1400
Ethyl Benzene	1.1	Not Detected	4.7	Not Detected
m,p-Xylene	1.1	Not Detected	4.7	Not Detected
o-Xylene	1.1	Not Detected	4.7	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	105	70-130
Toluene-d8	103	70-130
4-Bromofluorobenzene	91	70-130



Air Toxics

Client Sample ID: P2SVE-EFFC5-092215

Lab ID#: 1509395-03A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092812	Date of Collection:	9/22/15 9:30:00 AM
Dil. Factor:	2.46	Date of Analysis:	9/28/15 04:17 PM
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)
Vinyl Chloride	1.2	Not Detected	3.1
Chloroethane	4.9	Not Detected	13
1,1-Dichloroethene	1.2	1.2	4.9
Acetone	12	Not Detected	29
Methylene Chloride	12	Not Detected	43
trans-1,2-Dichloroethene	1.2	Not Detected	4.9
1,1-Dichloroethane	1.2	4.2	5.0
2-Butanone (Methyl Ethyl Ketone)	4.9	Not Detected	14
cis-1,2-Dichloroethene	1.2	2.4	4.9
Chloroform	1.2	Not Detected	6.0
1,1,1-Trichloroethane	1.2	150	6.7
Carbon Tetrachloride	1.2	Not Detected	7.7
Benzene	1.2	Not Detected	3.9
1,2-Dichloroethane	1.2	Not Detected	5.0
Trichloroethene	1.2	11	6.6
Toluene	1.2	Not Detected	4.6
1,1,2-Trichloroethane	1.2	Not Detected	6.7
Tetrachloroethene	1.2	47	8.3
Ethyl Benzene	1.2	Not Detected	5.3
m,p-Xylene	1.2	Not Detected	5.3
o-Xylene	1.2	Not Detected	5.3

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	98	70-130
Toluene-d8	103	70-130
4-Bromofluorobenzene	94	70-130



Air Toxics

Client Sample ID: P1SVE-EFFC2-092215

Lab ID#: 1509395-04A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092813	Date of Collection: 9/22/15 9:50:00 AM		
Dil. Factor:	2.26	Date of Analysis: 9/28/15 04:40 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.1	Not Detected	2.9	Not Detected
Chloroethane	4.5	Not Detected	12	Not Detected
1,1-Dichloroethene	1.1	3.4	4.5	14
Acetone	11	Not Detected	27	Not Detected
Methylene Chloride	11	Not Detected	39	Not Detected
trans-1,2-Dichloroethene	1.1	Not Detected	4.5	Not Detected
1,1-Dichloroethane	1.1	210	4.6	860
2-Butanone (Methyl Ethyl Ketone)	4.5	Not Detected	13	Not Detected
cis-1,2-Dichloroethene	1.1	9.6	4.5	38
Chloroform	1.1	Not Detected	5.5	Not Detected
1,1,1-Trichloroethane	1.1	450	6.2	2500
Carbon Tetrachloride	1.1	Not Detected	7.1	Not Detected
Benzene	1.1	Not Detected	3.6	Not Detected
1,2-Dichloroethane	1.1	Not Detected	4.6	Not Detected
Trichloroethene	1.1	30	6.1	160
Toluene	1.1	Not Detected	4.2	Not Detected
1,1,2-Trichloroethane	1.1	Not Detected	6.2	Not Detected
Tetrachloroethene	1.1	28	7.7	190
Ethyl Benzene	1.1	Not Detected	4.9	Not Detected
m,p-Xylene	1.1	Not Detected	4.9	Not Detected
o-Xylene	1.1	Not Detected	4.9	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	101	70-130
Toluene-d8	100	70-130
4-Bromofluorobenzene	94	70-130



Air Toxics

Client Sample ID: P1SVE-EFFC2-092215 DUP

Lab ID#: 1509395-05A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092814	Date of Collection:	9/22/15 9:50:00 AM	
Dil. Factor:	2.40	Date of Analysis:	9/28/15 05:16 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.2	Not Detected	3.1	Not Detected
Chloroethane	4.8	Not Detected	13	Not Detected
1,1-Dichloroethene	1.2	3.8	4.8	15
Acetone	12	Not Detected	28	Not Detected
Methylene Chloride	12	Not Detected	42	Not Detected
trans-1,2-Dichloroethene	1.2	Not Detected	4.8	Not Detected
1,1-Dichloroethane	1.2	210	4.8	840
2-Butanone (Methyl Ethyl Ketone)	4.8	Not Detected	14	Not Detected
cis-1,2-Dichloroethene	1.2	9.2	4.8	37
Chloroform	1.2	Not Detected	5.8	Not Detected
1,1,1-Trichloroethane	1.2	440	6.5	2400
Carbon Tetrachloride	1.2	Not Detected	7.6	Not Detected
Benzene	1.2	Not Detected	3.8	Not Detected
1,2-Dichloroethane	1.2	Not Detected	4.8	Not Detected
Trichloroethene	1.2	31	6.4	170
Toluene	1.2	Not Detected	4.5	Not Detected
1,1,2-Trichloroethane	1.2	Not Detected	6.5	Not Detected
Tetrachloroethene	1.2	19	8.1	130
Ethyl Benzene	1.2	Not Detected	5.2	Not Detected
m,p-Xylene	1.2	Not Detected	5.2	Not Detected
o-Xylene	1.2	Not Detected	5.2	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	98	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	92	70-130



Air Toxics

Client Sample ID: P1SVE-EFFC3-092215

Lab ID#: 1509395-06A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092817	Date of Collection: 9/22/15 10:25:00 AM		
Dil. Factor:	4.63	Date of Analysis: 9/28/15 06:51 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	2.3	Not Detected	5.9	Not Detected
Chloroethane	9.3	Not Detected	24	Not Detected
1,1-Dichloroethene	2.3	2.6	9.2	10
Acetone	23	Not Detected	55	Not Detected
Methylene Chloride	23	Not Detected	80	Not Detected
trans-1,2-Dichloroethene	2.3	Not Detected	9.2	Not Detected
1,1-Dichloroethane	2.3	11	9.4	43
2-Butanone (Methyl Ethyl Ketone)	9.3	Not Detected	27	Not Detected
cis-1,2-Dichloroethene	2.3	10	9.2	41
Chloroform	2.3	Not Detected	11	Not Detected
1,1,1-Trichloroethane	2.3	530	13	2900
Carbon Tetrachloride	2.3	Not Detected	14	Not Detected
Benzene	2.3	Not Detected	7.4	Not Detected
1,2-Dichloroethane	2.3	Not Detected	9.4	Not Detected
Trichloroethene	2.3	14	12	73
Toluene	2.3	Not Detected	8.7	Not Detected
1,1,2-Trichloroethane	2.3	Not Detected	13	Not Detected
Tetrachloroethene	2.3	18	16	120
Ethyl Benzene	2.3	Not Detected	10	Not Detected
m,p-Xylene	2.3	Not Detected	10	Not Detected
o-Xylene	2.3	Not Detected	10	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	99	70-130
Toluene-d8	101	70-130
4-Bromofluorobenzene	91	70-130



Air Toxics

Client Sample ID: FIELD BLANK-092215

Lab ID#: 1509395-07A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092816	Date of Collection:	9/22/15 10:30:00 AM	
Dil. Factor:	2.18	Date of Analysis:	9/28/15 06:16 PM	
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	1.1	Not Detected	2.8	Not Detected
Chloroethane	4.4	Not Detected	12	Not Detected
1,1-Dichloroethene	1.1	Not Detected	4.3	Not Detected
Acetone	11	Not Detected	26	Not Detected
Methylene Chloride	11	Not Detected	38	Not Detected
trans-1,2-Dichloroethene	1.1	Not Detected	4.3	Not Detected
1,1-Dichloroethane	1.1	Not Detected	4.4	Not Detected
2-Butanone (Methyl Ethyl Ketone)	4.4	Not Detected	13	Not Detected
cis-1,2-Dichloroethene	1.1	Not Detected	4.3	Not Detected
Chloroform	1.1	Not Detected	5.3	Not Detected
1,1,1-Trichloroethane	1.1	1.2	5.9	6.7
Carbon Tetrachloride	1.1	Not Detected	6.8	Not Detected
Benzene	1.1	Not Detected	3.5	Not Detected
1,2-Dichloroethane	1.1	Not Detected	4.4	Not Detected
Trichloroethene	1.1	Not Detected	5.8	Not Detected
Toluene	1.1	Not Detected	4.1	Not Detected
1,1,2-Trichloroethane	1.1	Not Detected	5.9	Not Detected
Tetrachloroethene	1.1	Not Detected	7.4	Not Detected
Ethyl Benzene	1.1	Not Detected	4.7	Not Detected
m,p-Xylene	1.1	Not Detected	4.7	Not Detected
o-Xylene	1.1	Not Detected	4.7	Not Detected

Container Type: 1 Liter Summa Canister

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	101	70-130
Toluene-d8	103	70-130
4-Bromofluorobenzene	91	70-130



Air Toxics

Client Sample ID: Lab Blank

Lab ID#: 1509395-08A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092807	Date of Collection: NA		
Dil. Factor:	1.00	Date of Analysis: 9/28/15 01:23 PM		
Compound	Rpt. Limit (ppbv)	Amount (ppbv)	Rpt. Limit (ug/m3)	Amount (ug/m3)
Vinyl Chloride	0.50	Not Detected	1.3	Not Detected
Chloroethane	2.0	Not Detected	5.3	Not Detected
1,1-Dichloroethene	0.50	Not Detected	2.0	Not Detected
Acetone	5.0	Not Detected	12	Not Detected
Methylene Chloride	5.0	Not Detected	17	Not Detected
trans-1,2-Dichloroethene	0.50	Not Detected	2.0	Not Detected
1,1-Dichloroethane	0.50	Not Detected	2.0	Not Detected
2-Butanone (Methyl Ethyl Ketone)	2.0	Not Detected	5.9	Not Detected
cis-1,2-Dichloroethene	0.50	Not Detected	2.0	Not Detected
Chloroform	0.50	Not Detected	2.4	Not Detected
1,1,1-Trichloroethane	0.50	Not Detected	2.7	Not Detected
Carbon Tetrachloride	0.50	Not Detected	3.1	Not Detected
Benzene	0.50	Not Detected	1.6	Not Detected
1,2-Dichloroethane	0.50	Not Detected	2.0	Not Detected
Trichloroethene	0.50	Not Detected	2.7	Not Detected
Toluene	0.50	Not Detected	1.9	Not Detected
1,1,2-Trichloroethane	0.50	Not Detected	2.7	Not Detected
Tetrachloroethene	0.50	Not Detected	3.4	Not Detected
Ethyl Benzene	0.50	Not Detected	2.2	Not Detected
m,p-Xylene	0.50	Not Detected	2.2	Not Detected
o-Xylene	0.50	Not Detected	2.2	Not Detected

Container Type: NA - Not Applicable

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	99	70-130
Toluene-d8	102	70-130
4-Bromofluorobenzene	94	70-130



Air Toxics

Client Sample ID: CCV

Lab ID#: 1509395-09A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092802	Date of Collection:	NA
Dil. Factor:	1.00	Date of Analysis:	9/28/15 09:56 AM

Compound	%Recovery
Vinyl Chloride	99
Chloroethane	100
1,1-Dichloroethene	97
Acetone	93
Methylene Chloride	103
trans-1,2-Dichloroethene	101
1,1-Dichloroethane	103
2-Butanone (Methyl Ethyl Ketone)	102
cis-1,2-Dichloroethene	100
Chloroform	98
1,1,1-Trichloroethane	99
Carbon Tetrachloride	99
Benzene	99
1,2-Dichloroethane	100
Trichloroethene	101
Toluene	101
1,1,2-Trichloroethane	99
Tetrachloroethene	94
Ethyl Benzene	104
m,p-Xylene	108
o-Xylene	108

Container Type: NA - Not Applicable

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	103	70-130
Toluene-d8	103	70-130
4-Bromofluorobenzene	98	70-130



Air Toxics

Client Sample ID: LCS

Lab ID#: 1509395-10A

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092803	Date of Collection:	NA
Dil. Factor:	1.00	Date of Analysis:	9/28/15 10:33 AM
Compound	%Recovery	Method	Limits
Vinyl Chloride	90	70-130	
Chloroethane	92	70-130	
1,1-Dichloroethene	85	70-130	
Acetone	82	70-130	
Methylene Chloride	90	70-130	
trans-1,2-Dichloroethene	79	70-130	
1,1-Dichloroethane	91	70-130	
2-Butanone (Methyl Ethyl Ketone)	90	70-130	
cis-1,2-Dichloroethene	98	70-130	
Chloroform	88	70-130	
1,1,1-Trichloroethane	87	70-130	
Carbon Tetrachloride	88	70-130	
Benzene	86	70-130	
1,2-Dichloroethane	87	70-130	
Trichloroethene	94	70-130	
Toluene	86	70-130	
1,1,2-Trichloroethane	84	70-130	
Tetrachloroethene	79	70-130	
Ethyl Benzene	89	70-130	
m,p-Xylene	91	70-130	
o-Xylene	92	70-130	

Container Type: NA - Not Applicable

Surrogates	%Recovery	Method	Limits
1,2-Dichloroethane-d4	103	70-130	
Toluene-d8	103	70-130	
4-Bromofluorobenzene	95	70-130	



Air Toxics

Client Sample ID: LCSD

Lab ID#: 1509395-10AA

MODIFIED EPA METHOD TO-14A GC/MS FULL SCAN

File Name:	17092804	Date of Collection:	NA
Dil. Factor:	1.00	Date of Analysis:	9/28/15 10:56 AM

Compound	%Recovery	Method Limits
Vinyl Chloride	91	70-130
Chloroethane	95	70-130
1,1-Dichloroethene	88	70-130
Acetone	87	70-130
Methylene Chloride	91	70-130
trans-1,2-Dichloroethene	78	70-130
1,1-Dichloroethane	92	70-130
2-Butanone (Methyl Ethyl Ketone)	89	70-130
cis-1,2-Dichloroethene	97	70-130
Chloroform	88	70-130
1,1,1-Trichloroethane	87	70-130
Carbon Tetrachloride	89	70-130
Benzene	88	70-130
1,2-Dichloroethane	87	70-130
Trichloroethene	94	70-130
Toluene	88	70-130
1,1,2-Trichloroethane	87	70-130
Tetrachloroethene	82	70-130
Ethyl Benzene	90	70-130
m,p-Xylene	92	70-130
o-Xylene	92	70-130

Container Type: NA - Not Applicable

Surrogates	%Recovery	Method Limits
1,2-Dichloroethane-d4	102	70-130
Toluene-d8	104	70-130
4-Bromofluorobenzene	95	70-130

C



C



Appendix C

**Third Quarter 2015
Phase1/Phase 2 AS/SVE
System Operations Data
Sheets**

DAILY DOCUMENTATION SHEET

Control Panel Touch Screen

DAILY DOCUMENTATION SHEET

Control Panel Touch Screen

DATE	7-4-15	7-5-15	7-6-15	7-7-15	7-8-15	7-9-15	7-10-15	7-11-15	7-12-15	7-13-15
TIME	9:30 AM	11:15 AM	8:02 AM	8:20 AM	2:00 PM	7:57 AM	10:10 AM	7:40 AM	10:20 AM	1:53 PM
OBSERVER'S INITIALS	DL	JL	RK	RL	PA	KK	RK	DE	KA	DE

ALARMS

HOURS METERS

B-701 SVE (hrs)	36413	36439	36459	36483	36512	36531	36557	36579	36605	36633
C-2201 SPRG (hrs)	31046	31072	31093	31118	31147	31166	31192	31214	31241	31269
F-2501 H-XCH (hrs)	31046	31072	31093	31118	31147	31166	31192	31214	31241	31269

ANALOGS

MV-701 SVE POS (%)	26	26	26	26	26	26	26	26	26	26	26
PT-701 SVE (-wc)	-81.6	-81.6	81.6	81.6	81.5	68.0	81.6	81.6	68.0	81.6	81.6
PT-2501 SPRG (psi)	13.3	11.9	12.9	12.9	12.9	12.7	12.9	13.0	12.8	12.9	12.9

SET POINTS

SET POINTS 2

SET POINTS 3

SET POINTS 4

MV-701 SVE POS (%) 30 30 30 30 30 30 30 30 30 30 30

DAILY DOCUMENTATION SHEET

Control Panel Touch Screen

DATE	6-24-15	6-25-15	6-26-15	6-27-15	6-28-15	6-29-15	6-30-15	7-1-15	7-2-15	7-3-15
TIME	1:31 P.M.	8:22 A.M.	8:23 A.M.	9:13 A.M.	10:00 A.M.	0845	8:59 A.M.	855	8:19 A.M.	8:25 A.M.
OBSERVER'S INITIALS	DE	PK	PK	DF	AC	A.H.	PK	ND	PK	K4

ALARMS

Shut Down Alarm Code	NA	NA	NA	NA	NA	N/A	NA	N/A	NA	NA
Non-critical Alarm Code	NA	NA	NA	NA	NA	N/A	NA	N/A	NA	NA

HOURS METERS

B-701 SVE (hrs)	36177	36196	36220	36245	36269	36292	36316	36340	36364	36388
C-2201 SPRG (hrs)	30808	30827	30852	30877	30902	30925	30949	30973	30997	31021
F-2501 H-XCH (hrs)	30808	30827	30852	30877	30902	30925	30949	30973	30977	31021

ANALOGS

MV-701 SVE POS (%)	26	26	26	26	26	26	26	26	26	26	26
PT-701 SVE (-wc)	81.6	81.6	81.6	81.6	-81.6	-81.6	81.6	81.6	81.6	81.6	68.0
PT-2501 SPRG (psi)	12.9	12.9	13.1	13.0	12.9	16.7	12.9	13.0	12.9	12.6	

SET POINTS

SET POINTS-2

SET POINTS 3

SVON-2801 SPRG (min)	60	60	60	60	60	60	60	60	60	60
SVON-2802 SPRG (min)	60	60	60	60	60	60	60	60	40	60
SVON-2803 SPRG (min)	180	180	180	180	180	180	180	180	180	180
SPRG DELAY (min)	0	0	0	0	15	0	0	0	0	0

SET POINTS A

MV-701 SVE POS (%) 30 30 30 30 38 30 31 30 30 30 30

WEEKLY DOCUMENTATION SHEET
SYSTEM COMPONENTS

DATE	6-8-15	6-15-15	6-22-15	6-29-15	7-6-15	7-14-15	9-22-15	9-28-15	10-5-15
TIME	8:43 AM	2:40 PM	8:16 AM	0845	8:04 AM	8:36 AM	1255	7:30 AM	8:11 AM
OBSERVER'S INITIALS	TK	KA	TK	A.LL	TK	TK	NP	TK	TK

SYSTEM LEAKS, EXCESSIVE OR UNFAMILIAR NOISE, MOISTURE, ETC.

Comments and Notes	NA	NA	NA	N/A	NA	NA	N/A	NA	NA
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SOIL VAPOR EXTRACTION (SVE)

Hours of Operation (hrs)	35790	35964	36024	36292	36459	36651	36702	36844	37012
Inlet Vacuum (-wc)	88	88	89	90	90	89	84	92	92
Pre-Filter Vacuum (-wc)	76.2	76	75.7	76.0	78.9	78.5	77.5	78.2	80.3
Post-Filter Vacuum (-wc)	78	78	78	76	80	80	70	79	80
Outlet Pressure (wc)	7	7	7	6.0	6.0	6.0	7.0	6.0	6.0
Outlet Temperature (°F)	160	164	162	160	175	170	150	164	158
Outlet Magnehelic* (in H ₂ O)	0.9	0.9	0.9	0.9	0.9	0.9	0.9	0.9	0.9
Water Level Sight Glass (in)	0	0	0	0	0	0	0	0	0

AIR SPARGE (SPRG)

Hours Operation (hrs)	30418	30593	30755	30915	31093	31287	31339	31483	31652
Oil Sight Glass (half pt.)	OK								

HEAT EXCHANGER (H-XCH)

Hours Operation (hrs)	30418	30593	30755	30925	3093	31287	31339	31483	31652
Inlet Temperature (°F)	205	216	205	200	210	215	205	210	200
Inlet Pressure (psi)	16	16	15.5	16.0	17.0	16	12	16	16.5
Outlet Temperature (°F)	110	112	110	105	129	130	112	109	108
Outlet Pressure (psi)	13.5	13	12	12.5	14.0	135	14.0	13.5	14
Outlet Magnehelic* (in H ₂ O)	4.0+	4.0+	4.0+	4.0+	4.+	4+	4+	4+	4+

ELECTRICAL USAGE (see display panel below main breaker and next to control panel)

Kilowatts (kwh)	17323	174700	175543	176809	177305	178320	178919	179668	180563
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* Keen plastic pinch valves on tubing to magnehelic gauge closed except when taking a reading.

MONTHLY DOCUMENTATION SHEET
SVE MANIFOLD

DATE	11/14/14	1-9-15	1-29-15	2-25-15	3-13-15	5-15-15	6-29-15	7-16-15	9/22/15
TIME	1055	0955	0850	1000	1045	0945	0845	0835	1255
INITIALS	NP/AS	B.H.	A.H.	A.H.	A.H.	W.H.	A.H.	A.H.	NP

/ MAGNEHELIC GAUGE*

CELL 1	SVE-1 (in H ₂ O)	1.1	1.1	1.1	1.1	1.0	1.0	1.2	1.2	1.2
	SVE-2 (in H ₂ O)	1.5	1.6	1.5	1.5	1.4	1.5	1.5	+0.5 _{0.5}	1.5
CELL 2	SVE-3 (in H ₂ O)	1.2	1.3	1.2	1.3	1.2	1.2	1.2	1.2	1.2
	SVE-4 (in H ₂ O)	1.4	1.4	1.3	1.4	1.6	1.2	1.3	1.2	1.2
CELL 3	SVE-5 (in H ₂ O)	1.1	1.2	1.0	0.9	1.0	0.9	1.2	1.2	1.1
	SVE-6 (in H ₂ O)	0.7	1.0	0.9	0.3	0.5	0.6	1.0	1.0	1.0

VACUUM GAUGE

CELL 1	SVE-1 (-wc)	30	30	30	30	32	28	32	33	32
	SVE-2 (-wc)	30	28	30	30	30	26	30	30	30
CELL 2	SVE-3 (-wc)	30	29	30	30	30	28	30	30	29
	SVE-4 (-wc)	30	29	18	16	20	18	30	30	29
CELL 3	SVE-5 (-wc)	36	36	36	36	36	36	36	37	36
	SVE-6 (-wc)	40	40	36	38	38	38	38	38	38

* Keep plastic pinch valves on tubing to magnehelic gauge closed except when taking a reading.

MONTHLY DOCUMENTATION SHEET
AIR SPARGE MANIFOLD

DATE	11/14/14	11/9/15	11/24/15	12/25/15	1/1/16	1/15/15	6/28/15	7/16/15	9/22/15	
TIME	10:53	09:55	08:50	10:00	10:45	09:45	08:45	08:35	12:00	
INITIALS	NMAS	R.H.	A.H.	R.H.	R.H.	R.H.	R.H.	R.H.	NP	
ROTOMETER										
CELL 1	AS-1 (scfm)	17	21	20	19	20	20	20	20	
	AS-2 (scfm)	20.5	21	21	21	21	20	21	21	
	AS-3 (scfm)	21	21	20	20.5	20	20	20	20	
	AS-4 (scfm)	19	19	20	19	20	20	19	20	
	AS-5 (scfm)	18	20	20	19	20	20	20	20	
CELL 2	AS-6 (scfm)	16	16	16	16	17	16	16	16	
	AS-7 (scfm)	16	16	16	16	17	16	16	16	
	AS-8 (scfm)	16	16	16	16	16	16	16	16	
	AS-9 (scfm)	16	16	16	16	16	16	16	16	
	AS-10 (scfm)	16.5	16.5	16.5	16.5	17	16.5	16.5	16.5	
CELL 3	AS-11 (scfm)	17	16	20	18	21	16	16	17	18
	AS-12 (scfm)	16	16	24	23.5	24	19	19	19	18
	AS-13 (scfm)	15.5	16	22	21	22	17.5	18	17	16.5
	AS-14 (scfm)	21.5	23	30	26	27.5	22.5	22.5	22	21.5
	AS-15 (scfm)	20.5	21.5	Not working	18.5	10.5	23.0	23.0	23	23
PRESSURE GAUGE										
CELL 1	AS-1 (psi)	11.5	11	10.0	10.5	9.5	10	10	10.5	10.5
	AS-2 (psi)	11	11	10.0	10.5	10	10	10	10.5	10.5
	AS-3 (psi)	10	10.5	9.5	10	9.5	10	10	10.0	10.0
	AS-4 (psi)	11	11	10.0	10.5	10	10	10	10.5	10.5
	AS-5 (psi)	11.5	11.5	10.0	10.5	10	10.5	10	10.5	10.5
CELL 2	AS-6 (psi)	11	11	10.0	10.5	10	10	10	10.5	10.5
	AS-7 (psi)	11	11	10.0	10.5	10	10	10	10.5	10.5
	AS-8 (psi)	11	11	10.0	10.5	10	10	10	10.5	10.5
	AS-9 (psi)	11	11	10.0	10.5	10	10	10	10.5	10.5
	AS-10 (psi)	11.5	11.5	10.0	10.5	10	10.5	10	10.5	10.5
CELL 3	AS-11 (psi)	11.5	11.5	9.5	10	9.5	10	9.5	10.5	10
	AS-12 (psi)	13.5	13.5	10.5	14	11	13.5	13.5	13.5	13.5
	AS-13 (psi)	13	13	10.0	10	11	13.5	11.5	13.0	12.0
	AS-14 (psi)	13.5	14	10.5	12.5	11	12.5	13.5	12.0	12.0
	AS-15 (psi)	13	13.5	13.0	13.5	13.5	13.5	13.5	13.5	12.5

MONTHLY DOCUMENTATION SHEET
WELL HEAD GAUGES

DATE	3-13-15	5-15-15	6-29-15	7-16-15	9/22/15					
TIME	1045	0945	0845	0835	1255					
INITIALS	A.H.	A.H.	A.H.	A.H.	NP					

CELL 1	AS-1 (psi)	Covered	8.75	8.8	8.9	9.36				
	AS-2 (psi)	8.0	8.0	8.0	8.0	9.05				
	AS-3 (psi)	4.0	4.0	4.5	4.5	9.18				
	AS-4 (psi)	5.0	2.0	2.0	2.0	9.24				
	AS-5 (psi)	12.0	12.0	12.0	12.0	9.19				
CELL 2	AS-6 (psi)	8.0	8.0	8.0	8.0	9.98				
	AS-7 (psi)	8.0	8.0	8.0	8.0	9.73				
	AS-8 (psi)					9.24				
	AS-9 (psi)					9.24				
	AS-10 (psi)					9.24				
CELL 3	AS-11 (psi)	8.0	6.0	5.0	9.0	8.84				
	AS-12 (psi)	10.5	10.5	10.0	10.5	9.54				
	AS-13 (psi)	7.0	4.0	2.5	7.0	9.56				
	AS-14 (psi)	10.5	10.5	10	10.5	9.48				
	AS-15 (psi)	Broken	10	9.53	9.95	9.55				

CELL 1	SVE-1 (in H ₂ O)	6.5	6.5	6.5	6.5	0.01				
	SVE-2 (in H ₂ O)	2.0	2.0	1.0	1.5	0.01				
CELL 2	SVE-3 (in H ₂ O)									
	SVE-4 (in H ₂ O)									
CELL 3	SVE-5 (in H ₂ O)	2.5	10	10	2.5	0.34				
	SVE-6 (in H ₂ O)	1.0	8	8.1	8.3	0.27				

DAILY DOCUMENTATION SHEET

BUILDINGS 1 - 2

Control Panel Touch Screen

DATE	9-24-15	09-25-15	9-26-15	9-27-15	9-28-15	9-29-15	9-30-15	10-1-15	10-2-15
TIME	11:11	9:07 AM	5:39 AM	8:06 AM	7:44 AM	9:06 AM	8:56 AM	7:06 AM	
OBSERVER'S INITIALS	AA	KA	PK	PK	PK	PK	PK	PK	PL

WEATHER CONDITIONS

Indoor Room Temp (°F), Outdoor Conditions: (Rain, Snow, Clear, Overcast, etc.)	80°F Sunny	76°F SUNNY	80°F DARK	76°F Cloudy - RAIN	74°F RAIN	66°F CLEAR	66°F CLEAR	62°F CLEAR
	S	S	DARK	Cloudy - RAIN	RAIN	CLEAR	CLEAR	CLEAR

ALARMS

Alarm Code	T	F	Name	NONE	NONE	N/A	N/A	N/A	N/A

P&ID

PDT-701 SVE (-wc)	M	0.06	0.13	0.13	0.29	0.25	0.14	0.14	0.11
PT-701 SVE (-wc)		-49	-51	-58	-61	-63	-59	-59	-58
PT-702 SVE (-wc)		60.1	59.9	58.5	57.4	57.6	57.8	58.0	59.0
PT-2201 SPRG (psi)	D	17.3	17.1	16.7	16.4	16.4	16.7	17.2	12.5
P-401 PUMP (cycles)	O	49	49	49	49	49	49	49	49

P&ID2

PDT-801 SVE (-wc)	W	0.29	0.31	0.28	0.32	0.32	0.34	0.35	0.38
PT-801 SVE (-wc)	A	-125	-135	-141	-144	-146	-139	-137	-140
PT-802 SVE (-wc)	A	49.2	47.8	47.1	46.8	46.9	48.4	47.8	48.0
PT-2301 SPRG (psi)		0	0	0.0	0.0	0.0	0.0	0.0	0.0
P-501 PUMP (cycles)		225	225	227	229	232	235	238	241

DAILY DOCUMENTATION SHEET

BUILDINGS 1 - 2

Control Panel Touch Screen

DATE	9-24-15	9-25-15	9-26-15	9-27-15	9-28-15	9-29-15	9-30-15	10-1-15	10-2-15
TIME	11:11	9:07 AM	5:39 AM	8:06 AM	7:44 AM	9:06 AM	8:52 AM	7:06 AM	
OBSERVER'S INITIALS	A.IH	KA	TR	TR	PK	TR	TR	TR	BL
HOURS METERS	S								
B-701 SVE (hrs)	X	6571	65.93	6613	6640	6643	6689	6713	6735
C-2201 SPRG (hrs)	X	6504	6525	6546	6572	6596	6621	6645	6667
C-2202 SPRG (hrs)	S	6198	6220	6241	6267	6291	6316	6340	6362
B-801 SVE (hrs)	T	6507	6529	6550	6576	6600	6625	6649	6671
C-2301 SPRG (hrs)	T	6489	6511	6532	6558	6582	6607	6631	6653
C-2302 SPRG (hrs)	I	6489	6515	6532	6558	6582	6607	6631	6653
SET POINTS									
PAL-701 SVE (wc)	11	-27	-27	-27	-27	-27	-27	-27	-27
PAH-702 SVE (wc)		100	100.0	100.0	100.0	100.0	100.0	100.0	100.0
PAL-702 SVE (wc)	D	10	10.0	10.0	10.0	10.0	10.0	10.0	10.0
PAH-2201 SPRG (psi)	D	30	30	30.0	30.0	30.0	30.0	30.0	30.0
PAL-2201 SPRG (psi)	O	0	0.0	0.0	0.0	0.0	0.0	0.0	0.0
PAL-801 SVE (wc)	O	-25	-25	-25	-25	-25	-25	-25	-25
PAH-802 SVE (wc)	W	100	100.0	100.0	100.0	100.0	100.0	100.0	100.0
PAL-802 SVE (wc)	O	10	10.0	10.0	10.0	10.0	10.0	10.0	10.0
PAH-2301 SPRG (psi)	N	30	30	30.0	30.0	30.0	30.0	30.0	30.0
PAL-2301 SPRG (psi)	O	0	0.0	0.0	0.0	0.0	0.0	0.0	0.0
SV-2801 SPRG (min)		40	40	40	40	40	40	40	40
SV-2802 SPRG (min)		20	20	20	20	20	20	20	20
SV-2901 SPRG (min)		20	20	20	20	20	20	20	20
SV-2902 SPRG (min)		20	20	20	20	20	20	20	20

DAILY DOCUMENTATION SHEET

BUILDINGS 1 - 2

Control Panel Touch Screen

DATE	7-9-15	7-10-15	7-11-15	7-12-15	7-13-15	7-14-15	7-15-15	7-16-15	9/22/15	9/23-15
TIME	8:03 AM	1145	7:48 AM	10:27 AM	2:00 PM	8:34 AM	1030		1100	
OBSERVER'S INITIALS	RL	A.H.	DE	KA	DE	RL	A.H.	NP		

WEATHER CONDITIONS

Indoor Room Temp (°F), Outdoor Conditions: (Rain, Snow, Clear, Overcast, etc.)	CLEAR	clear	Cloudy	Cloudy	CLEAR	CLEAR	cloudy	CLEAR	S Y S
	74°F	84°F	76°F	80°F	90°F	74°F	76°F	78°	T C M

ALARMS

Alarm Code	NA	n/a	NA	NA	NA	NA	NA	NA	T C M
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P&ID

PDT-701 SVE (-wc)	0.08	0.06	0.10	0.12	0.07	0.09	0.10	0.07	J
PT-701 SVE (-wc)	-64	-57	-60	-63	-70	-64	-55	-49	J
PT-702 SVE (-wg)	55.6	57.2	56.9	55.0	52.7	55.4	55.1	61.4	O
PT-2201 SPRG (psi)	14.0	13.2	13.0	12.9	10.6	11.6	14.2	15.5	O
P-401 PUMP (cycles)	49	49	49	49	49	49	49	49	W

P&ID2

PDT-801 SVE (-wc)	0.27	0.29	0.32	0.30	0.27	0.25	0.28	0.25	
PT-801 SVE (-wc)	-180	-138	-132	-135	-145	-139	-136	-122	J
PT-802 SVE (-wc)	47.8	46.7	48.2	47.4	46.7	47.8	47.3	49.3	
PT-2301 SPRG (psi)	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0.0	
P-501 PUMP (cycles)	219	219	220	220	220	220	220	223	

Transmitter not recording.

DAILY DOCUMENTATION SHEET

BUILDINGS 1 - 2

Control Panel Touch Screen

DATE	7-9-15	7-10-15	7-11-15	7-12-15	7-13-15	7-15-15	7-16-15	7-22-15	7-23-15
TIME	8:03 AM	1145	7:48 AM	15:27 AM	2:00 P.M.	8:24 AM	1030	1100	
OBSERVER'S INITIALS	RK	A.14	DE	KA	DE	RK	A.4	NP	

HOURS METERS

B-701 SVE (hrs)	6361	6387	6403	6423	6451	6485	6504	6506	S
C-2201 SPRG (hrs)	6294	6320	6336	6356	6383	6418	6436	6439	Y
C-2202 SPRG (hrs)	6131 (044)	6131	6131	6131	6131	6131	6131	6134	S
B-801 SVE (hrs)	6295	6323	6338	6359	6386	6421	6439	6442	S
C-2301 SPRG (hrs)	6277	6305	6321	6341	6368	6403	6421	6424	T
C-2302 SPRG (hrs)	6277	6305	6320	6341	6368	6403	6421	6424	C

SET POINTS

DAILY DOCUMENTATION SHEET

BUILDINGS 1 - 2

Control Panel Touch Screen

DATE	6-30-15	7-1-15	7-2-15	7-3-15	7-4-15	7-5-15	7-6-15	7-7-15	7-8-15
TIME	8:58AM	0853	1230	8:15	9:35	11:00AM	8:10AM	-	1330
OBSERVER'S INITIALS	RK	A. H.	A. H.	KA	OL	OL	RK	A. H.	A. H.

WEATHER CONDITIONS

Indoor Room Temp (°F), Outdoor Conditions: (Rain, Snow, Clear, Overcast, etc.)	CLEAR	clear	clear	SUNNY	SUNNY	SUNNY	partly cloudy	System down	Partly Cloudy
	70°	60°F	70°F	72°F	70°	78°	81°F		79°F

ALARMS

Alarm Code	NA	N/A	N/A	NA	NA	NA	NA		N/A
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P&ID

PDT-701 SVE (-wc)	0.12	0.10	0.05	0.07	0.08	0.09	0.09		0.10
PT-701 SVE (-wc)	-65	-58	-53	-60	-60	-67	-70		-58
PT-702 SVE (-wc)	55.9	58.9	59.3	58.9	54.4	53.6	53.5		56.9
PT-2201 SPRG (psi)	13.2	14.8	14.5	13.9	10.8	13.4	13.5		14.9
P-401 PUMP (cycles)	49	49	49	49	49	49	49		49

P&ID2

PDT-801 SVE (-wc)	0.29	0.32	0.29	0.23	0.29	0.29	0.31		0.29
PT-801 SVE (-wc)	-144	-127	-124	-135	-141	-143	-148		-129
PT-802 SVE (-wc)	47.3	49.7	49.6	48.6	47.0	45.0	45.8		47.7
PT-2301 SPRG (psi)	0.0	0.0	0.0	0.0	0.0	0.0	0.0		0.0
P-501 PUMP (cycles)	218	218	218	218	219	219	219		219

Transmitter not recording.

DAILY DOCUMENTATION SHEET

BUILDINGS 1 - 2

Control Panel Touch Screen

DATE	6-30-15	7-1-15	7-2-15	7-3-15	7-4-15	7-5-15	7-6-15	7-7-15	7-8-15
TIME	8:36AM	0853	1230	8:16AM	9:36AM	11:00AM	8:19AM	-	1330
OBSERVER'S INITIALS	TK	IAH	PA	KA	DL	DL	PC	PA	PA
HOURS METERS									
B-701 SVE (hrs)	6222	6223	6246	6266	6291	6317	6338	S	6343
C-2201 SPRG (hrs)	6155	6156	6179	6199	6224	6250	6271	X	6275
C-2202 SPRG (hrs)	6130	6131	6131	6131	6131	6131	6131	S	6131
B-801 SVE (hrs)	6154	6157	6180	6200	6225	6251	6272	T	6276
C-2301 SPRG (hrs)	6138	6139	6162	6182	6207	6233	6274	G	6259
C-2302 SPRG (hrs)	6138	6139	6162	6182	6207	6233	6274	M	6258
SET POINTS									
PAL-701 SVE (wc)	-27	-27	-27	-27	-27	-27	-27		-27
PAH-702 SVE (wc)	100.0	100.0	100.0	100.0	100.0	100.0	100.0	D	100.0
PAL-702 SVE (wc)	10.0	10.0	10.0	10.0	10.0	10.0	10.0	O	10.0
PAH-2201 SPRG (psi)	30.0	30.0	30.0	30.0	30.0	30.0	30.0	6	30.0
PAL-2201 SPRG (psi)	0.0	0.0	0.0	0.0	0.0	0.0	0.0	0	0.0
PAL-801 SVE (wc)	-25	-25	-25	-25	-25	-25	-25	N	-25
PAH-802 SVE (wc)	100.0	100.0	100.0	100.0	100.0	100.0	100.0		100.0
PAL-802 SVE (wc)	10.0	10.0	10.0	10.0	10.0	10.0	10.0		10.0
PAH-2301 SPRG (psi)	30.0	30.0	30.0	30.0	30.0	30.0	30.0		30.0
PAL-2301 SPRG (psi)	0.0	0.0	0.0	0.0	0.0	0.0	0.0		0.0
SV-2801 SPRG (min)	40	40	40	400	400	40.0	40		40
SV-2802 SPRG (min)	20	20	20	20	20	20	20	V	20
SV-2901 SPRG (min)	20	20	20	20	20	20	20		20
SV-2902 SPRG (min)	20	20	20	20	30	20	20		20

WEEKLY DOCUMENTATION SHEET
BUILDING 1
SYSTEM COMPONENTS

9/22/15

DATE	6-15-15	6-22-15	6-29-15	7-6-15	7-13-15	7-16-15 ¹⁰⁰	9-28-15	10-5-15
TIME	2:15PM	8:31AM	0930	8:16AM	8:25AM	1030 ¹⁴	8:03AM	8:23AM
OBSERVER'S INITIALS	KH	RY	A.H	RL	RL	RP A.H. & RH	RK	PK

SYSTEM LEAKS, EXCESSIVE OR UNFAMILIAR NOISE, MOISTURE, ETC.

Comments and Notes	NO	NA	N/A	NA	NA	N/A	NA	NA
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SOIL VAPOR EXTRACTION (SVE) BLOWER B-701

Pre-Filter Vacuum (-wc)	55	75	52	60	55	40	50	51
Post-Filter Vacuum (-wc)	90	90	84	94	88	70	83	84
Inlet Magnehelic* (in H ₂ O)	.3	.2	0.2	0.2	0.2	0.15	0.15	0.2
Inlet Vacuum (-wc)	80	-82	82	-82	Gauge Break	Gauge Break	Gauge Break	Gauge Break

SOIL VAPOR EXTRACTION (SVE) HEAT EXCHANGER (H-XCH)

Inlet Pressure (wc)	22	22	23	18	20	28	40	20
Inlet Temperature (°F)	140	102	125	130	125	125	125	115
Outlet Pressure (wc)	9	8	5.5	6	8	8	9	8
Outlet Temperature (°F)	109	102	94	95	90	98	92	74
Water Level Sight Glass (in)	0	0	0	0	8"	0	Ø	Ø

AIR SPARGE (SPRG) COMPRESSOR C-2201

Upper Oil Sight Glass (half pt.)	OK							
Lower Oil Sight Glass (half pt.)	OK							

AIR SPARGE (SPRG) HEAT EXCHANGER (H-XCH)

Inlet Pressure (psi)	25	15	15	14.5	160	2	17.5	13
Inlet Temperature (°F)	250	255	195	200	189	135	205	190
Outlet Pressure (psi)	24	14.5	15.5	15	15	2	18.0	12.5
Outlet Temperature (°F)	122	118	82	86	80	95	98	90

* Keep plastic pinch valves on tubing to magnehelic gauge closed except when taking a reading.

WEEKLY DOCUMENTATION SHEET
BUILDING 2
SYSTEM COMPONENTS

DATE	6-29-15	7-6-15	7-14-15	9/22/15	9-22-15	10-5-15		
TIME	0905	8:10AM	8:34AM	1040	7:58AM	8:18AM		
OBSERVER'S INITIALS	A. H	FL	FL	NP	TR	RK		

SYSTEM LEAKS, EXCESSIVE OR UNFAMILIAR NOISE, MOISTURE, ETC.

Comments	NA	NA	NA	NA	NA	NA		
and Notes	74°F	80°F	82°F	78°F	79°F	66°F		

SOIL VAPOR EXTRACTION (SVE) BLOWER B-801

Pre-Filter Vacuum (-wc)	58	65	55	50	60	62		
Post-Filter Vacuum (-wc)	60	70	80	55	66	68		
Inlet Magnehelic* (in H ₂ O)	0.40	0.40	03.5	0.30	0.4	0.405		
Inlet Vacuum (-wc)	60	-40	-60	-60	-60	-60		

SOIL VAPOR EXTRACTION (SVE) HEAT EXCHANGER (H-XCH)

Inlet Pressure (wc)	48	46	48	52	47	48		
Inlet Temperature (°F)	125	125	130	125	126	124		
Outlet Pressure (wc)	1.0	1	1	2	1	2		
Outlet Temperature (°F)	96	98	100	100	98	86		
Water Level Sight Glass (in)	8 1/2	0	9"	0	24"	20"		

AIR SPARGE (SPRG) COMPRESSOR C-2301

Upper Oil Sight Glass (half pt.)	OK	OK	OK	OK	OK	OK		
Lower Oil Sight Glass (half pt.)	OK	OK	OK	OK	OK	OK		

AIR SPARGE (SPRG) HEAT EXCHANGER (H-XCH)

Inlet Pressure (psi)	18	17.5	16.5	18.0	17.5	16		
Inlet Temperature (°F)	210	212	220	215	215	198		
Outlet Pressure (psi)	16.5	16.5	16.5	16.5	16.0	15.5		
Outlet Temperature (°F)	118	118	126	120	119	108		

* Keep plastic pinch valves on tubing to magnehelic gauge closed except when taking a reading.

MONTHLY DOCUMENTATION SHEET
BUILDING 1
SVE MANIFOLD

DATE	6-24-15	7/16-15	9/22/15								
TIME	0930	1030	1100								
INITIALS	A.H.	A.H.	NP								

MAGNEHELIC GAUGE*

SVE-15 (in H ₂ O)	0.0	0.0	0.0								
SVE-14 (in H ₂ O)	0.0	0.0	0.0								
SVE-13 (in H ₂ O)	0.0	0.0	0.0								
SVE-11 (in H ₂ O)	0.25	0.40	0.25								
SVE-9 (in H ₂ O)	0.55	0.40	0.40								

VACUUM GAUGE

SVE-15 (-wc)	50	55	40								
SVE-14 (-wc)	53	55	40								
SVE-13 (-wc)	55	60	40								
SVE-11 (-wc)	18	20	14.5								
SVE-9 (-wc)	50	55	35								

* plastic pinch valves on tubing to magnehelic gauge closed except when taking a reading.

MONTHLY DOCUMENTATION SHEET
BUILDING 1 UPPER CELL
AIR SPARGE MANIFOLD

DATE	5-15-15	6-29-15	7-16-15	9/22/15					
TIME	0905	0930	1030	1100					
INITIALS	A-H.	A-H.	A-H.	NP					

ROTOMETER

UPPER CELL	AS-47 (scfm)	0	0	10	16				
	AS-38 (scfm)	10	5	5	6				
	AS-37 (scfm)	16	8	7	11				
	AS-23 (scfm)	15	7.5	6	11				
	AS-24 (scfm)	14	7.0	6.5	10				
	AS-39 (scfm)	16	8.0	7	11				
	AS-48 (scfm)	17.5	9.0	8.5	13				
	AS-52 (scfm)	15	7.5	7	10.5				
	AS-56 (scfm)	14	7.5	6.5	10				
	AS-54 (scfm)	22	10	10	14.5				
	AS-58 (scfm)	12.5	6.5	5.5	9				

PRESSURE GAUGE

UPPER CELL	AS-47 (psi)	11.5	0	12	17				
	AS-38 (psi)	17	13	14	14				
	AS-37 (psi)	8.5	7.5	7	8				
	AS-23 (psi)	13	11	10.5	12.5				
	AS-24 (psi)	13	12	11	13				
	AS-39 (psi)	12	11	10	12				
	AS-48 (psi)	11.5	11	10	11.5				
	AS-52 (psi)	11	10.5	9.5	11				
	AS-56 (psi)	11	10.0	9.5	11				
	AS-54 (psi)	14	13.0	14	13.5				
	AS-58 (psi)	9.5	12.0	11.5	10				

*Note: 6-29-15 reading low due to C-2207 not in operation,
 7-16-15 11

MONTHLY DOCUMENTATION SHEET
BUILDING 1 LOWER CELL
AIR SPARGE MANIFOLD

DATE	5-15-15	6-29-15	7-16-15	9/22/15					
TIME	0905	0930	1030	1100					
INITIALS	A.H.	A.H.	A.H.	NP					

ROTOMETER

LOWER CELL	AS-49 (scfm)	15	8	7.5	11				
	AS-44 (scfm)	14	8	7.5	12				
	AS-31 (scfm)	12	7	7.5	10				
	AS-32 (scfm)	12	7.5	7.5	10				
	AS-45 (scfm)	14.5	6	6.5	11				
	AS-51 (scfm)	15	7	7.5	11.5				
	AS-55 (scfm)	18	10	10	14.5				
	AS-53 (scfm)	17	8	7.5	12				
	AS-59 (scfm)	14.5	8	7.5	11				
	AS-57 (scfm)	16.5	6	6	11.5				
	AS-50 (scfm)	22	10	9.5	17				

PRESSURE GAUGE

LOWER CELL	AS-49 (psi)	10.5	10	10	10.5				
	AS-44 (psi)	12.5	10.5	11	12				
	AS-31 (psi)	10	10	10	10				
	AS-32 (psi)	9	10	10	9.5				
	AS-45 (psi)	11.5	10.5	11	11.5				
	AS-51 (psi)	10.5	10	10	10.5				
	AS-55 (psi)	10.5	10	10	10.5				
	AS-53 (psi)	11.5	10	10.5	11.5				
	AS-59 (psi)	11	10	10.5	11				
	AS-57 (psi)	12	10.5	11	12				
	AS-50 (psi)	12	10.5	11	12				

4-18-15 reading low due to C-2202 not in operation
7-16-15 11

Dan n-a

MONTHLY DOCUMENTATION SHEET
BUILDING 1
PID MEASUREMENTS

DATE	1-4-15	2-25-15	2-13-15	5-25-15	(6-24-15)	7-16-15	9/22/15					
TIME	1255	0930	0945	0905	0930	1030	1100					
INITIALS	A.H.	H.H.	A.H.	A.H.	A.H.	A.H.	NP					

SVE Effluent	0.9	0.4	1.9	0.0	0.0	0.0	0.0					
SVE - 15	moisture	moisture	moisture	0.0	0.0	moisture	0.0					
SVE - 14				0.0	moisture		0.0					
SVE - 13				0.0	0.0		0.0					
SVE - 11				0.0	0.0		0.0					
SVE - 9	↓	↓	↓	0.0	0.0	↓	0.0					
Exhaust	0.9	0.4	1.9	0.0	0.0	0.0	0.0					

MONTHLY DOCUMENTATION SHEET
BUILDING 2
SVE MANIFOLD

DATE	6-24-15	7-16-15	9/22/15							
TIME	0905	1020	1040							
INITIALS	A. H.	A. H.	N.P.							

MAGNEHELIC GAUGE*

SVE-10 (in H ₂ O)	0.35	0.25	0.25							
SVE-12 (in H ₂ O)	0.60	0.60	0.50							
SVE-8 (in H ₂ O)	0.60	0.70	0.50							
SVE-7 (in H ₂ O)	0.40	0.35	0.25							

VACUUM GAUGE

SVE-10 (-wc)	55	80	90							
SVE-12 (-wc)	55	55	48							
SVE-8 (-wc)	57	55	50							
SVE-7 (-wc)	65	68	65							

ELECTRICAL USAGE

Kilowatts (kwh)	1338.00	3165.40	37182.70							
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* Keep plastic pinch valves on tubing to magnehelic gauge closed except when taking a reading.

MONTHLY DOCUMENTATION SHEET
BUILDING 2 UPPER CELL
AIR SPARGE MANIFOLD

DATE	5-15-15	10-26-15	7-16-15	9/22/15					
TIME	1026	0905	1020	1440					
INITIALS	A.H.	A.H.	A.H.	NP					

ROTOMETER

UPPER CELL	AS-20 (scfm)	15	15	15					
	AS-26 (scfm)	9	8	8.5	9				
	AS-16 (scfm)	17	16	17	17				
	AS-18 (scfm)	11	10	10	10				
	AS-22 (scfm)	13.5	13	13	13.5				
	AS-28 (scfm)	13	12	13	13.5				
	AS-30 (scfm)	20	18	19	19.5				
	AS-36 (scfm)	16	16	15.5	16				
	AS-42 (scfm)	15.5	16	10	10.5				
	AS-40 (scfm)	16	15.5	19	19.5				
	AS-34 (scfm)	17	27	27	29				

PRESSURE GAUGE

UPPER CELL	AS-20 (psi)	10	10	10	10				
	AS-26 (psi)	10	10.5	10	10.5				
	AS-16 (psi)	10	10	10	10				
	AS-18 (psi)	14.5	14.5	15	15				
	AS-22 (psi)	10	10	10	10				
	AS-28 (psi)	10.5	10.5	10.5	10.5				
	AS-30 (psi)	10	10	10	10.5				
	AS-36 (psi)	11.5	11.0	11.5	12				
	AS-42 (psi)	10.5	10.5	10	10.5				
	AS-40 (psi)	10	10.5	10.5	11				
	AS-34 (psi)	10	11.0	11.5	11.5				

MONTHLY DOCUMENTATION SHEET
BUILDING 2 LOWER CELL
AIR SPARGE MANIFOLD

DATE	5-15-15	6-24-15	7-16-15	9/22/15			
TIME	1030	0905	1020	1140			
INITIALS	A.H.	A.H.	A.H.	NP			

ROTOMETER

LOWER CELL	AS-27 (scfm)	14	14	15	15.5			
	AS-25 (scfm)	15.5	15	17	18			
	AS-17 (scfm)	15	14	15	16			
	AS-19 (scfm)	15	15	16	17			
	AS-21 (scfm)	18	20	20	22			
	AS-29 (scfm)	10.5	20	15	16.5			
	AS-43 (scfm)	16.5	10	11	18.5			
	AS-46 (scfm)	21	21	22	24			
	AS-41 (scfm)	13	12	12.5	13			
	AS-33 (scfm)	17	17	18	19.5			
	AS-35 (scfm)	13.5	13	13.5	14.5			

PRESSURE GAUGE

LOWER CELL	AS-27 (psi)	12	11.5	10.5	11			
	AS-25 (psi)	11	10.5	10.5	10.5			
	AS-17 (psi)	11	10.5	11.0	11.5			
	AS-19 (psi)	13	12.5	13	12.5			
	AS-21 (psi)	12.5	11.5	12	11.5			
	AS-29 (psi)	14	12.5	13.5	13.5			
	AS-43 (psi)	11	10.5	10.5	10.5			
	AS-46 (psi)	12.5	12	12	12			
	AS-41 (psi)	10.5	10	10.5	10			
	AS-33 (psi)	11.5	11	11	11			
	AS-35 (psi)	11.5	11	11	10.5			

MONTHLY DOCUMENTATION SHEET
BUILDING 2
PID MEASUREMENTS

DATE	1-9-15	2-25-15	3-13-15	5-15-15	6-29-15	7-16-15	9/22/15					
TIME	1240	0950	1000	1030	0905	1020	1040					
INITIALS	A.H.	A.H.	A.H.	A.H.	A.H.	A.H.	NP					

SVE Effluent	0.00	0.00	0.90	0.00	0.00	0.0	0.00					
SVE - 10	moisture	moisture	moisture	0.00	moisture	moisture	0.0					
SVE - 12				0.00	moisture		0.0					
SVE - 8				0.00	moisture		0.0					
SVE - 7	↓	↓	↓	0.00	moisture	↓	0.0					
Exhaust	0.00	0.00	0.90	0.00	0.00	0.0	0.0					

D



D



Appendix D

Third Quarter 2015 Groundwater Sampling Data Sheets

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-4-15 Time: Start 1235 (24hr)
Project No: 60339110-4214 Finish 1345
Site Location: Rockford, Illinois
Weather: Sunny 70-80°F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 47.73 Screen interval(ft): 15 Approx. depth of pump intake(ft): 40
Water table depth (ft): 34.02 Casing type/diameter: 2" PVC Minimum purge volume: 6,70 (gals)
Water column length (ft): 13.71 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	3512-2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	1240		

(continued on back)

Sample Collector(s):

Date: 8-4-15

SAMPLE COLLECTION DATA

Well ID:

GMZ-01

Page 2 of 2

VOCS - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-5-15 Time: Start 1540 (24hr)
Project No: 60339110-4214 Finish 1645
Site Location: Rockford, Illinois
Weather: Sunny 75-85°F Collector(s): A. 1b/1g/1z

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 44.80 Screen interval(ft): 15 Approx. depth of pump intake(ft): 37
Water table depth (ft): 31.50 Casing type/diameter: 2" PVC Minimum purge volume: (6,10) (gals)
Water column length (ft): 1244 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotté	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	1540		

Time (24hr)	Purge Vol. (ml)	Temp. (°C)	pH	Spec. Cond. (µS/cm)	ORP (mV)	DO (mg/L)	Turbidity (NTU)	Flow Rate (ml/min)	Drawdown (feet)	Color/Odor
15:50	5000	16.55	7.67	1098	221.4	9.36	14.5	500	31.56	clear/inone
15:55	7500	16.79	7.67	1017	193.0	8.50	7.47	500	31.65	
16:00	10,000	16.60	7.61	1015	1718.0	8.29	4.69	500	31.65	
16:05	12,500	16.63	7.61	1014	163.1	8.29	3.78	500	31.65	
16:10	15,000	16.41	7.61	1011	154.6	8.23	3.33	500	31.65	
16:15	17,500	16.60	7.61	1009	147.2	8.18	2.99	500	31.65	
16:20	20,000	16.32	7.61	1009	141.8	8.26	2.49	500	31.65	
16:25	22,500	16.40	7.61	1008	140.7	8.23	1.98	500	31.65	
16:30	25,000	16.43	7.61	1010	136.2	8.20	1.88	500	31.65	

(continued on back)

Sample Collector(s):

allotted

Date: 8-5-15

SAMPLE COLLECTION DATA

Well ID:

GMZ-02

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VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-5-15 Time: Start 1405 (24hr)
Project No: 60339110-4214 Finish 1515
Site Location: Rockford, Illinois
Weather: Sunny 75-85°F Collector(s): A. Helgatis

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 44.60 Screen interval(ft): 15 Approx. depth of pump intake(ft): 37
Water table depth (ft): 30.95 Casing type/diameter: 2" PVC Minimum purge volume: (6.70) (gals)
Water column length (ft): 13.65 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	<u>1405</u>		

Time (24hr)	Purge Vol. (ml)	Temp. (°C)	pH	Spec. Cond. (µS/cm)	ORP (mV)	DO (mg/L)	Turbidity (NTU)	Flow Rate (ml/min)	Drawdown (feet)	Color/Odor
1405	5000	17.18	7.88	1016	123.7	8.69	20.8	500	30.95	clear/none
1420	7500	17.27	7.68	1019	113.9	7.52	9.87	500	31.05	
1425	10,000	17.13	7.73	1020	108.0	7.37	7.15	500	31.05	
1430	12,500	16.88	7.77	1022	105.8	7.32	5.71	500	31.05	
1435	15,000	16.85	7.75	1019	107.0	7.24	4.12	500	31.05	
1440	17,500	16.82	7.79	1019	106.7	7.28	3.61	500	31.05	
1445	20,000	17.02	7.74	1016	111.3	7.19	2.88	500	31.05	
1450	22,500	16.82	7.81	1020	115.3	7.24	2.74	500	31.05	
1455	25,000	16.90	7.81	1019	120.0	7.22	2.23	500	31.05	
1500	17,500	17.01	7.81	1019	123.9	7.19	2.13	500	31.05	

* Turbidity not stable, removed 3 turb. volumes, collect sample.

(continued on back)

Sample Collector(s):

Date: 8-5-15

SAMPLE COLLECTION DATA

Well ID:

GMZ-03

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VOCs - Volatile organic compounds

G - Glass

HCl - Hydrchloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility

Date: 8-5-15

Time: Start
Finish

0950 (24hr)
105

Project No: 60339110-4214

Finish

Site Location: Rockford, Illinois

Weather: Sunny 70-80°F

Collector(s): A. Hellmuth

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 45.30 Screen interval(ft): 15 Approx. depth of pump intake(ft): 38
Water table depth (ft): 29.27 Casing type/diameter: 2" PVC Minimum purge volume: 7.90 (gals)
Water column length (ft): 16.03
(calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	0955		

Time (24hr)	Purge Vol. (ml)	Temp. (°C)	pH	Spec. Cond. (µS/cm)	ORP (mV)	DO (mg/L)	Turbidity (NTU)	Flow Rate (ml/min)	Drawdown (feet)	Color/Odor
1005	5000	17.18	8.09	1666.2	147	8.17	80.2	500	29.27	cloudy/none
1010	7500	17.30	8.06	1666.7	134.9	7.97	55.6	500	29.27	↓
1015	10,000	17.12	8.09	1651.0	124.9	7.86	33.1	500	29.27	clear/none
1020	12,500	17.02	8.12	1658.8	106.5	7.64	15.2	500	29.30	
1025	15,000	17.06	8.14	165.2	110.5	7.36	9.66	500	29.30	
1030	17,500	17.10	8.16	1648.5	106.6	7.19	6.65	500	29.30	
1035	20,000	17.02	8.10	1647.7	103.5	7.13	5.10	500	29.30	
1040	22,500	17.19	8.17	1644.4	101.5	7.12	4.47	500	29.30	
1045	25,000	17.04	8.17	1645.5	104.6	7.13	3.36	500	29.30	
1050	17,500	16.95	8.17	1649.6	94.4	7.24	3.04	500	29.30	
1055	30,000	16.93	8.18	1647.7	97.7	7.25	2.93	500	29.30	↓

(continued on back)

Sample Collector(s):

Allen West

Date: 8-5-15

SAMPLE COLLECTION DATA

Well ID:

GMZ-04

Page 2 of 2

VOCS - Volatile organic compounds

G - Glass

HCl - Hydrchloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-5-15 Time: Start 0730 (24hr)
Project No: 60339110-4214 Finish 0840
Site Location: Rockford, Illinois
Weather: Sunny 66° - 70°F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 46.79 Screen interval(ft): 15 Approx. depth of pump intake(ft): 39
 Water table depth (ft): 29.33 Casing type/diameter: 4" ~~2"~~ SS Minimum purge volume: 34.20 (gals)
 Water column length (ft): 17.46 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F666
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	044-0702
Begin purge at	0735		

(continued on back)

Sample Collector(s):

Date: 8-5-15

SAMPLE COLLECTION DATA

Well ID:

MW-07FGA

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VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-4-15 Time: Start 1545 (24hr)
Project No: 6039110-4214 Finish 1710
Site Location: Rockford, Illinois
Weather: Sunny 70-80°F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 47.50 Screen interval(ft): 15 Approx. depth of pump intake(ft): 40
Water table depth (ft): 30.12 Casing type/diameter: 2" SS Minimum purge volume: 8.50 (gals)
Water column length (ft): 17.38 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	04F66
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	<u>1550</u>		

Time (24hr)	Purge Vol. (ml)	Temp. (°C)	pH	Spec. Cond. (µS/cm)	ORP (mV)	DO (mg/L)	Turbidity (NTU)	Flow Rate (ml/min)	Drawdown (feet)	Color/Odor
1600	5000	18.21	7.03	1063	22.1	0.74	24.0	500	30.12	clear/none
1605	7500	18.19	6.94	1028	29.7	0.57	15.3	500	30.14	
1610	10,000	18.24	6.94	1056	22.8	0.64	10.05	500	30.14	
1615	13,500	18.29	6.95	1064	23.5	0.52	8.71	500	30.14	
1620	15,000	18.24	6.95	1071	27.3	0.40	6.38	500	30.14	
1625	17,500	18.25	6.916	1075	35.1	0.29	4.98	500	30.14	
1630	20,000	18.46	6.93	1078	41.0	0.24	4.58	500	30.14	
1635	22,500	18.70	6.96	1081	46.8	0.25	4.02	500	30.14	
1640	25,000	18.86	6.96	1085	51.9	0.23	3.68	500	30.14	
1645	27,500	18.19	6.916	1087	54.8	0.20	3.05	500	30.14	
1650	30,000	17.67	6.95	1090	55.5	0.18	2.03	500	30.14	
1655	30,500	17.98	6.95	1090	58.2	0.26	3.45	500	30.14	

(continued on back)

Sample Collector(s):

alterative

Date: 8-4-15

SAMPLE COLLECTION DATA

Well ID: MW-203

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VOCs - Volatile organic compounds

G · Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-6-15 Time: Start 0720 (24hr)
Project No: 6039110-4214 Finish 0830
Site Location: Rockford, Illinois
Weather: Sunny 60-70°F Collector(s): A. Hellas -

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 44.72 Screen interval(ft): 25 Approx. depth of pump intake(ft): 37
Water table depth (ft): 31.71 Casing type/diameter: 2" PVC Minimum purge volume: 6.4 (gals)
Water column length (ft): 13.01

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09FL66
	Lamotte	2020	2532-3512
	Lamotte	Smart 2 Colorimeter	0344-0702
Begin purge at	0730		

* Turbidity next stable, remove 3 well volumes, collect sample

(continued on back)

Sample Collector(s):

Allen Bent

Date: 8-6-15

SAMPLE COLLECTION DATA

Well ID:

PMW-01

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VOCs - Volatile organic compounds

G - Glass
HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-10-15 Time: Start 0840 (24hr)
Project No: 60339110-4214 Finish 0935
Site Location: Rockford, Illinois
Weather: Sunny 68°-70° F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 41.85 Screen interval(ft): 25 Approx. depth of pump intake(ft): 37
Water table depth (ft): 31.64 Casing type/diameter: 2" PVC Minimum purge volume: 50 (gals)
Water column length (ft): 10.21

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	0845		

(continued on back)

Sample Collector(s):

Date: 8-6-15

SAMPLE COLLECTION DATA

Well ID:

PMW-02

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-4-15 Time: Start 1350 (24hr)
Project No: 60339110-4214 Finish 1435
Site Location: Rockford, Illinois
Weather: Sunny 70-80°F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 39.50 Screen interval(ft): 15 Approx. depth of pump intake(ft): 32
 Water table depth (ft): 32.08 Casing type/diameter: 2" PVC Minimum purge volume: 3,60 (gals)
 Water column length (ft): 7.42 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	CGF66
	Lamotte	2020	2537-372
	Lamotte	Smart 2 Colorimeter	0244~0202
Begin purge at	1355		

(continued on back)

Sample Collector(s):

all over

Date: 8-4-15

SAMPLE COLLECTION DATA

Well ID:

SMW-01

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VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-4-15 Time: Start 1440 (24hr)
Project No: 60339110-4214 Finish 1540
Site Location: Rockford, Illinois
Weather: Sunny 70-80°F Collector(s): A. Holzatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 40.00 Screen interval(ft): 15 Approx. depth of pump intake(ft): 32
Water table depth (ft): 28.59 Casing type/diameter: 2" PVC Minimum purge volume: 5.60 (gals)
Water column length (ft): 11.41 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F666
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244 - 0702
Begin purge at	1445		

(continued on back)

Sample Collector(s):

all set

Date: 8-4-15

SAMPLE COLLECTION DATA

Well ID:

SMW-02

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass

HCl - Hydrchloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-16-15 Time: Start 0940 (24hr)
Project No: 60339110-4214 Finish 1040
Site Location: Rockford, Illinois
Weather: Sunny (65-75°F) Collector(s): A. Holkatt

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 42.50 Screen interval(ft): 15 Approx. depth of pump intake(ft): 35
Water table depth (ft): 31.22 Casing type/diameter: 2" PVC Minimum purge volume: 5.60 (gals)
Water column length (ft): 11.28 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	2537-3519
	Lamotte	Smart 2 Colorimeter	0246-5702
Begin purge at	0945		

at ORP, 10 outflow, larger flow volumes. Scum left collector.

(continued on back)

Sample Collector(s):

Date: 8-6-15

SAMPLE COLLECTION DATA

Well ID: **SMW-04**

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VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-4-15 Time: Start 1135 (24hr)
Project No: 60339110-4214 Finish 1230
Site Location: Rockford, Illinois
Weather: Sunny 70° 80°F Collector(s): A-Hole 12

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 41.85 Screen interval(ft): 15 Approx. depth of pump intake(ft): 34
Water table depth (ft): 31.42 Casing type/diameter: 2" PVC Minimum purge volume: 5,10 (gals)
Water column length (ft): 10.43 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	1244-0702
Begin purge at	1140		

(continued on back)

Sample Collector(s):

Allen Art

Date: 8-4-15

SAMPLE COLLECTION DATA

Well ID:

SMW-08

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-5-15 Time: Start 0845 (24hr)
Project No: 60339110-4214 Finish 0940
Site Location: Rockford, Illinois
Weather: Sunny 65-75°F Collector(s): A. Haller

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 41.16 Screen interval(ft): 15 Approx. depth of pump intake(ft): 35
Water table depth (ft): 30.31 Casing type/diameter: 2" SS Minimum purge volume: 5.30 (gals)
Water column length (ft): 10.85 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	69F66
	Lamotte	2020	8537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	0850		

(continued on back)

Sample Collector(s):

Alfred

Date: 8-5-15

SAMPLE COLLECTION DATA

Well ID:

SMW-19

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VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-5-15 Time: Start 1300 (24hr)
Project No: 60339110-4214 Finish 1400
Site Location: Rockford, Illinois
Weather: Sunny 75-85°F Collector(s): A. Hallalz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 40.15 Screen Interval(ft): 15 Approx. depth of pump intake(ft): 33
Water table depth (ft): 30.41 Casing type/diameter: 2" PVC Minimum purge volume: 4.80 (gals)
Water column length (ft): 9.74
(calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F706
	Lamotte	2020	2532-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	1305		

(continued on back)

Sample Collector(s):

Date: 8-5-15

SAMPLE COLLECTION DATA

Well ID: **SMW-20**

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VOCS - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: <u>UTAS Plants 1/2 Facility</u>	Date: <u>8-5-15</u>	Time:	Start <u>110</u>	(24hr)
Project No: <u>60339110-4214</u>		Finish	<u>1215</u>	
Site Location: <u>Rockford, Illinois</u>				
Weather: <u>Sunny 70-80°F</u>	Collector(s): <u>A. Hollatz</u>			

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 41.40 Screen interval(ft): 15 Approx. depth of pump intake(ft): 34
Water table depth (ft): 29.85 Casing type/diameter: 2" PVC Minimum purge volume: 5.70 (gals)
Water column length (ft): 11.58 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0344-0702
Begin purge at	1115		

(continued on back)

Sample Collector(s):

Allen M. St

Date: 8-5-15

SAMPLE COLLECTION DATA

Well ID:

SMW-21

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VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-6-15 Time: Start 1105 (24hr)
Project No: 60339110-4214 Finish 1220
Site Location: Rockford, Illinois
Weather: Sunny 70-80°F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 446.00 Screen interval(ft): 15 Approx. depth of pump intake(ft): 38
 Water table depth (ft): 31.60 Casing type/diameter: 2" PVC Minimum purge volume: 7,10 (gals)
 Water column length (ft): 14.40 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09 Flodo
	Lamotte	2020	2527-3512
	Lamotte	Smart 2 Colorimeter	0244-0707
Begin purge at	1115		

Turbid or not stable, removed 2 mL 110 volumes, collected sample.

(continued on back)

Sample Collector(s):

albicans

Date: 8-6-15

SAMPLE COLLECTION DATA

Well ID:

RAMW-01

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VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-6-15 Time: Start 1225 (24hr)
Project No: 60339110-4214 Finish 1340
Site Location: Rockford, Illinois
Weather: Sunny 75-85° F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 44.84 Screen Interval(ft): 15 Approx. depth of pump intake(ft): 37
Water table depth (ft): 31.48 Casing type/diameter: 2" PVC Minimum purge volume: 66 (gals)
Water column length (ft): 13.36 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09E606
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	1230		

* Turbid, try ORP not stable, Remove Blv's (volume)
Collect sample.

(continued on back)

Sample Collector(s):

Allen Melt

Date: 10-8-14 8-6-15
1914

SAMPLE COLLECTION DATA

Well ID:

RAMW-02

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass

HCl - Hydrchloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-6-15 Time: Start 1350 (24hr)
Project No: 60339110-4214 Finish 1500
Site Location: Rockford, Illinois
Weather: Partly cloudy 75-85°F Collector(s): A-Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 45.70 Screen interval(ft): 15 Approx. depth of pump intake(ft): 38
Water table depth (ft): 31.29 Casing type/diameter: 2" PVC Minimum purge volume: 6.90 (gals)
Water column length (ft): 14.01

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	0537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	1355		

Time (24hr)	Purge Vol. (ml)	Temp. (°C)	pH	Spec. Cond. (µS/cm)	ORP (mV)	DO (mg/L)	Turbidity (NTU)	Flow Rate (ml/min)	Drawdown (feet)	Color/Odor
1405	5000	19.42	7.29	847	144.6	3.83	3.46	500	31.29	clear/inone
1410	7,500	19.41	7.26	852	130.1	3.30	3.14	500	31.30	
1415	10,000	19.37	7.26	851	120.2	3.27	3.52	500	31.30	
1420	12,500	19.45	7.26	854	112.9	3.26	2.55	500	31.30	
1425	15,000	19.44	7.27	855	105.1	3.29	3.05	500	31.30	
1430	17,500	19.07	7.27	861	102.7	3.35	3.35	500	31.30	
1435	20,000	18.84	7.27	859	104.0	3.34	2.42	500	31.30	
1440	22,500	18.87	7.27	863	107.8	3.35	2.06	500	31.30	
1445	25,000	18.53	7.26	864	112.4	3.37	1.94	500	31.30	
1450	27,500	18.43	7.26	867	116.0	3.34	1.67	500	31.30	▼

(continued on back)

Sample Collector(s):

Date: 8-6-15

SAMPLE COLLECTION DATA

Well ID:

RAMW-03

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass

HCl - Hydrchloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-7-15 Time: Start 115 (24hr)
Project No: 60339110-4214 Finish 1205
Site Location: Rockford, Illinois
Weather: Sunny 70-80° F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 45.00 Screen interval(ft): 15 Approx. depth of pump intake(ft): 41
Water table depth (ft): 31.14 Casing type/diameter: 2" PVC Minimum purge volume: 10.8 (gals)
Water column length (ft): 13.86
(calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F666
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	1120		

(continued on back)

Sample Collector(s):

Date: 8-7-15

SAMPLE COLLECTION DATA

Well ID: RAMW-04

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass
HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-7-15 Time: Start 0955 (24hr)
Project No: 60339110-4214 Finish 1110
Site Location: Rockford, Illinois
Weather: Sunny 65°-75°F Collector(s): A. Hallal 2

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 43.85 Screen interval(ft): 15 Approx. depth of pump intake(ft): 36
Water table depth (ft): 30.02 Casing type/diameter: 2" PVC Minimum purge volume: 68 (gals)
Water column length (ft): 13.83
(calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	04F66
	Lamotte	2020	2537-3572
	Lamotte	Smart 2 Coloriméter	0444-0702
Begin purge at	1000		

Time (24hr)	Purge Vol. (ml)	Temp. (°C)	pH	Spec. Cond. (µs/cm)	ORP (mV)	DO (mg/L)	Turbidity (NTU)	Flow Rate (ml/min)	Drawdown (feet)	Color/Odor
10/10	5000	16.83	7.84	1445	1816.4	8.78	10.20	500	30.00	clear/none
10/15	7500	16.90	7.23	1407	172.2	8.45	5.76	500	30.00	
10/20	10,000	17.09	7.74	1386	157.7	8.32	4.99	500	30.00	
10/25	12,500	17.18	7.80	1321	148.1	8.16	4.23	500	30.00	
10/30	15,000	17.75	7.83	1356	143.6	7.73	3.52	500	30.00	
10/35	17,500	17.40	7.82	1350	142.3	7.95	3.05	500	30.00	
10/40	20,000	17.21	7.82	1344	141.2	7.84	2.104	500	30.00	
10/45	22,500	17.52	7.83	1333	141.6	7.75	2.08	500	30.00	
10/50	25,000	17.83	7.83	1327	142.8	7.35	1.82	500	30.00	
10/55	27,500	17.53	7.82	1322	143.3	7.55	1.166	500	30.00	

(continued on back)

Sample Collector(s):

Date: 8-7-15

SAMPLE COLLECTION DATA

Well ID:

RAMW-05

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass
HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-7-15 Time: Start 0850 (24hr)
Project No: 60339110-4214 Finish 0950
Site Location: Rockford, Illinois
Weather: Sunny 65-75°F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 44.27 Screen interval(ft): 15 Approx. depth of pump intake(ft): 37
 Water table depth (ft): 30.07 Casing type/diameter: 2" PVC Minimum purge volume: 70 (gals)
 Water column length (ft): 14.20 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	2537-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	08:55		

(continued on back)

Sample Collector(s):

Date: 8-7-15

SAMPLE COLLECTION DATA

Well ID:

RAMW-06

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass

HCl - Hydrchloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility
Project No.: 60339110-4214
Site Location: Rockford, Illinois
Weather: Sunny 60 - 70 °F

Date: 8-7-15

Time: Start
Finish

0715 (24hr)
0820

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 48.100 Screen interval(ft): 15 Approx. depth of pump intake(ft): 41
 Water table depth (ft): 34.53 Casing type/diameter: 2" PVC Minimum purge volume: 6.9 (gals)
 Water column length (ft): 14.67 (calculations on reverse)

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	09F66
	Lamotte	2020	2C37-3512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at	0720		

(continued on back)

Sample Collector(s):

Date: 8-7-15

SAMPLE COLLECTION DATA

Well ID:

RAMW-07

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Ground Water Sample Collection Record

Client: UTAS Plants 1/2 Facility Date: 8-16-15 Time: Start 15:10 (24hr)
Project No: 60339110-4214 Finish 16:25
Site Location: Rockford, Illinois
Weather: Sunny 75-85°F Collector(s): A. Hollatz

1. WELL and WATER LEVEL DATA: (measured from Top of Casing)

Total well length (ft): 44.29 Screen interval(ft): 15 Approx. depth of pump intake(ft): 37
Water table depth (ft): 30.64 Casing type/diameter: 2" PVC Minimum purge volume: (6.70) (gals)
Water column length (ft): 13.65

2. WELL PURGE DATA

Purge/Sample Method: Proactive SS Monsoon Pump

Well is stable when readings stabilize to +/- 10% over three (3) consecutive readings collected at 5-minute intervals. If three (3) well volumes have been removed, and the readings have not stabilized, a sample shall be collected.

Field Testing Equipment Used:	Make	Model	Serial Number(s)
	YSI	556 MPS	091-66
	Lamottc	2020	2537-2512
	Lamotte	Smart 2 Colorimeter	0244-0702
Begin purge at 1515 1545A.M.			

Time (24hr)	Purge Vol. (ml)	Temp. (°C)	pH	Spec. Cond. (µS/cm)	ORP (mV)	DO (mg/L)	Turbidity (NTU)	Flow Rate (ml/min)	Drawdown (feet)	Color/Odor
1525	5000	18.91	7.25	1645	-190.3	0.89	28.4	500	30.64	Clear/none
1530	7500	18.92	7.31	1545	-181.3	0.50	24.5	500	30.94	
1535	10,000	18.88	7.30	1447	-165.3	0.35	15.7	500	30.94	
1540	12,500	18.88	7.27	1398	-134.5	0.32	12.3	500	30.94	
1545	15,000	18.08	7.25	1376	-134.7	0.37	10.14	500	30.94	
1550	17,500	18.64	7.24	1359	-129.3	0.29	09.53	500	30.94	
1555	20,000	18.69	7.23	1362	-129.8	0.28	7.26	500	30.94	
1600	22,500	18.73	7.22	1366	-127.5	0.24	6.83	500	30.94	
1605	25,000	18.81	7.21	1372	-125.9	0.24	6.37	500	30.94	
1610	27,500	18.84	7.20	1383	-122.9	0.23	5.77	500	30.94	

(continued on back)

Sample Collector(s):

Date: 8-6-15

SAMPLE COLLECTION DATA

Well ID:

RAMW-08

Page 2 of 2

VOCs - Volatile organic compounds

G - Glass

HCl - Hydrochloric acid

Project Name: UTAS Plants 1/2 Facility

Project Number: 60339110

Date: 8-4-15

Calibration Form

Parameter	Instrument		Standard		Standard Value @ C	Ambient Temp. C	Initial Value	Adjusted Value	Initials & Time	Comments
	Manf/Model	Serial No.	Manf/Model	SN/Exp. Date						
pH 4.00	YSI 556 MPS	09F66	NCL Lab	B441502/86 1-26-17	4.00 @ 25C	28.85	4.01	4.00	A.H.1047	
			NCL Lab	B347140805 8.5-16		24.01	7.03	7.00	A.H.1045	Post Cal
pH 7.00	YSI	15E100303	NCL Lab	B5G14D811 8-11-16	10.00 @ 25C	23.95	10.01	10.00	A.H.1049	
pH 10.00			YSI	11-16		23.92	1047	1000	A.H.1043	Post Cal
Specific Cond.	YSI	145100415 9-16		1,000 uS/cm @ 25C	24.01	239.1	244.0	1051		Post Cal
ORP				241.mV @ 15 C		24.01	239.1	244.0	1051	Post Cal
DO		H2O Saturated Air	—	100% H2O Sat. Air	24.20	113.0	96.2	A.H.1041	BP =	
Turbidity	LaMotte 2020	2537- 3512	—	—	0 NTU	NA	0.00	0.00	A.H.1053	
						NA	—	—		Post Cal
			—	—	10 NTU	NA	10.04	10.01	L.H.1053	
						NA	—	—		Post Cal

BP = Barometric Pressure (mmHg)

Project Name: UTA's Plants 1/2 Facility

Project Number: 60339110

Date: 8-5-15

Calibration Form

Parameter	Instrument		Standard		Standard Value @ C	Ambient Temp. C	Initial Value	Adjusted Value	Initials & Time	Comments
	Manf/Model	Serial No.	Manf/Model	SN/Exp. Date						
pH 4.00	YSI 556MP3 09F66	NCL Lab B441502106 B441408045 B5014D911 YSI 15E100303 11-16 145100415 9-16 H2O Saturated Air	NCL	B441502106	4.00 @ 25C	23.80	4.06	4.00	A.H07010	
pH 7.00				1-26-17		23.70	7.06	7.00	A.H0704	
pH 10.00				8-5-16						Post Cal
Specific Cond.				10.00 @ 25C		23.85	10.03	10.00	A.H0708	
ORP				1,000 uS/cm @ 25C		23.45	1004	1000	A.H0702	
DO				244 mV @ 15 C		23.46	239.5	244.0	A.H0710	
Turbidity	Lamotte 2020	2537- 3512	—	—	0 NTU	NA	0.02	0.00	A.H0658	
					10 NTU	NA	9.93	10.01	A.H0658	Post Cal
						NA				Post Cal

BP = Barometric Pressure (mmHg)

Project Name: UTA 5 ~~Plants~~ 1/2 Feb 1/11

Project Number: 60339110

Date: 8-5-15

Calibration Form

Parameter	Instrument		Standard		Standard Value @ C	Ambient Temp. C	Initial Value	Adjusted Value	Initials & Time	Comments
	Manf/Model	Serial No.	Manf/Model	SN/Exp. Date						
pH 4.00	YSI 556 MPS	09F66	NCL Lab	D441503134 1-26-17	4.00 @ 25C	25.40	4.07	4.00	A-H1520	
				B47140805 8-5-16		25.24	7.00	7.00	A-H1524	Post Cal
pH 7.00					7.00 @ 25C					Post Cal
pH 10.00			YSI	B5014D811 9-11-16	10.00 @ 25C	25.43	10.04	10.01	A-H1528	
Specific Cond.			YSI	15E100303 11-16	1,000 uS/cm @ 25C	25.18	1009	1000	1522A4	
ORP			YSI	145100415 9-16	244 mV @ 15 C	25.56	235.5	244.0	A-H1530	
DO			H2O Saturated Air	—	100% H2O Sat. Air	25.05	85.4	96.2	A-H1520	BP =
Turbidity	Lamotte 2020	2537 3512	—	—	0 NTU	NA	0.00	0.00	A-H1530	
					10 NTU	NA	9.93	10.01	A-H1530	Post Cal
					NA			—		Post Cal

BP = Barometric Pressure (mmHg)

Project Name: WTPS Plant 1/2 Facility

Project Number: 60339110

Date: 8-16-15

Calibration Form

Parameter	Instrument		Standard		Standard Value @ C	Ambient Temp. C	Initial Value	Adjusted Value	Initials & Time	Comments
	Manf/Model	Serial No.	Manf/Model	SN/Exp. Date						
pH 4.00	YSI 556 MPS	09F66	NCL Lab	B4411502126 1-26-17	4.00 @ 25C	24.62	4.00	4.00	A.140700	
pH 7.00				B471140805 8-5-16		24.76	7.01	7.00	A.140658	Post Cal
pH 10.00			↓	B5011410811 8-14-16	10.00 @ 25C	24.49	10.04	10.01	A.140702	Post Cal
Specific Cond.			YSI	15E100303 11-16		24.64	10.35	1000	A.140656	Post Cal
ORP			↓	143100415 9-16	244 mV @ 15°C	24.10	23.1	244.0	A.140704	Post Cal
DO			H2O Saturated Air	—		23.20	99.9	96.2	A.140654	BP =
Turbidity	LaMotte 2020	2537 3512	—		0 NTU	NA	0.02	0.00	A.140706	
					10 NTU	NA	10.08	10.00	A.140706	Post Cal
						NA				Post Cal

BP = Barometric Pressure (mmHg)

Project Name: UTA S. Plants 1/2 Facility

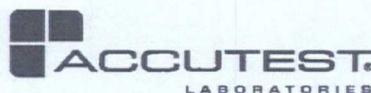
Project Number: 6033910

Date: 8-7-15

Calibration Form

Parameter	Instrument		Standard		Standard Value @ C	Ambient Temp. C	Initial Value	Adjusted Value	Initials & Time	Comments
	Manf/Model	Serial No.	Manf/Model	SN/Exp. Date						
pH 4.00	YSI 556 MPS	09F66	NCL Lab	B-141502126 1-26-17	4.00 @ 25C	24.77	4.03	4.00	AH 0701	Post Cal
pH 7.00				B47140805 8-5-16		24.84	6.96	7.00	AH 0659	
pH 10.00	YSI		↓	B501410811 8-11-16	10.00 @ 25C	24.72	10.05	10.01	AH 0703	Post Cal
Specific Cond.			YSI	1SE100203 11-16		24.62	1013	1000	AH 0657	
ORP	YSI		↓	143100415 9-16	244mv @ 15 C	24.30	246.8	244.0	AH 0705	Post Cal
DO			H2O Saturated Air	—		33.35	95.5	96.1	AH 0655	
Turbidity	LaMotte	2537 3512	—	—	0 NTU	NA	0.00	0.00	AB 0705	BP = Post Cal, BP = Post Cal
						NA	—	—	—	
			—	—	10 NTU	NA	9.87	10.01	AB 0705	
						NA	—	—	—	

BP = Barometric Pressure (mmHg)



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Client / Reporting Information		Project Information				Requested Analysis (see TEST CODE sheet)								Matrix Codes					
Company Name AECOM		Project Name: UTAS Plants H2 Facilities												DW - Drinking Water GW - Ground Water WW - Water SW - Surface Water SO - Soil SL - Sludge SED - Sediment OI - Oil LIQ - Other Liquid AIR - Air SOL - Other Solid WP - Wipe FB - Field Blank EB - Equipment Blank RB - Rinse Blank TB - Trip Blank					
Street Address 27755 Dierhard Suite 100		Street Rockford		Billing Information (if different from Report to)															
City Warrenville IL	State IL	Zip 60555	City Rockford	State IL	Company Name														
Project Contact Peter Hollatz	E-mail Peter.Hollatz@aecom.com	Project # 60339110	Street Address																
Phone # (630) 836-5386	Fax # (630) 836-1711	Client Purchase Order # 5948 ACW1	City Rockford				State IL	Zip 60555											
Sampler(s) Name(s) Allan Hollatz	Phone # (630) 836-1700	Project Manager Peter Hollatz	Attention:																
Accutest Sample #	Field ID / Point of Collection	MEOH/DI Vial #	Collection			Matrix	# of bottles	Number of preserved Bottles								VOC	LAB USE ONLY		
			Date 8-6-15	Time 1210	Sampled by AH			HCl	NaOH	HNO3	H2SO4	NONE	DI Water	MEOH	EDTA			CH3CO	
HSER-RAMW01-080615			8-6-15	1210	AH	GW	3	X								X			
HSER-FBLK02-080615			8-6-15	1105			3		X							X			
HSER-RAMW02-080615			8-6-15	1320			3	X								X			
HSER-MSD02-080615			8-6-15	1320			3		X							X			
HSER-RAMW02-080615			8-6-15	1320			3	X								X			
HSER-RAMW03-080615			8-6-15	1450			3	X								X			
HSER-RAMW08-080615			8-6-15	1610			3	X								X			
HSER-DUP02-080615			8-6-15	1600			3	X								X			
HSER-RAMW07-080715			8-7-15	0810			3	X								X			
HSER-RAMW06-080715			8-7-15	0935			3	X								X			
HSER-RAMW05-080715			8-7-15	1055			3	X								X			
HSER-RAMW04-080715			8-7-15	1155			3	X								X			
Turnaround Time (Business days)										Data Deliverable Information								Comments / Special Instructions	
<input type="checkbox"/> Std. 10 Business Days <input type="checkbox"/> 5 Day RUSH <input type="checkbox"/> 3 Day RUSH <input type="checkbox"/> 2 Day RUSH <input type="checkbox"/> 1 Day RUSH <input type="checkbox"/> other _____										Approved By (Accutest PM): / Date: <hr/> <hr/> <hr/> <hr/> <hr/> <hr/>								<input type="checkbox"/> Commercial "A" (Level 1) <input type="checkbox"/> Commercial "B" (Level 2) <input type="checkbox"/> FULLT1 (Level 3+4) <input type="checkbox"/> NJ Reduced <input type="checkbox"/> Commercial "C" <input type="checkbox"/> NYASP Category A <input type="checkbox"/> NYASP Category B <input type="checkbox"/> State Forms <input type="checkbox"/> EDD Format <input type="checkbox"/> Other _____ <input type="checkbox"/> NJ Data of Known Quality Protocol Reporting Commercial "A" = Results Only, Commercial "B" = Results + QC Summary NJ Reduced = Results + QC Summary + Partial Raw data	
Emergency & Rush T/A data available VIA Lablink										List of 13 VOC LEVEL 1 VOC Data									
Sample Custody must be documented below each time samples change possession, including courier delivery.																			
Relinquished by Sampler: Allen Hollatz		Date Time: 8-7-15 1800	Received By: 1	Relinquished By: 2				Date Time: 8-7-15 1800				Received By: 2							
Relinquished by Sampler: 3		Date Time:	Received By: 3	Relinquished By: 4				Date Time:				Received By: 4							
Relinquished by Sampler: 5		Date Time:	Received By: 5	Custody Seal #				<input type="checkbox"/> Intact	Preserved where applicable				On Ice	Comp.	<input type="checkbox"/>				



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Appendix E

Fourth Quarter 2015 Progress Report



Submitted to:
UTC Aerospace Systems
Rockford, IL

Submitted by:
AECOM
Warrenville, IL
60339110
December 2015

QUARTERLY PROGRESS REPORT – Fourth Quarter 2015 (September 2015 – November 2015)

UTC Aerospace Systems Plants 1/2 Facility
Southeast Rockford Groundwater Contamination
Superfund Site
2421 11th Street
Rockford, Illinois 61104
ILD981000417, ILD010219665
AECOM Project No. 60339110

This Quarterly Progress Report has been prepared on behalf of UTC Aerospace Systems (UTAS, fka Hamilton Sundstrand or HS) by AECOM Technical Services Inc. (AECOM). This report summarizes activities that occurred during the months of September, October, and November of the Fourth Quarter of 2015 at the above-referenced facility.

Progress Report- Reporting Quarters			
Q1	December	January	February
Q2	March	April	May
Q3	June	July	August
Q4	September	October	November

This report is the twenty-ninth in the series of Quarterly Progress Reports and consistent with United States Environmental Protection Agency (USEPA) approval of combining project reporting documents from a letter dated April 15, 2011. Quarterly Progress Reports are included as attachments to the Groundwater Management Zone (GMZ) Monitoring and System Performances Reports.

This Quarterly Progress Report follows the requirements outlined in Section X of the Consent Decree (CD) and includes the following:

Actions taken during the prior quarter to maintain compliance with the CD:

- Summaries of sampling results and tests.
- An identification of work plans and other deliverables completed in accordance with the CD.
- Actions scheduled for the next quarter.
- Information on the progress, percentage of completion, delays, and efforts to mitigate delays.
- Modifications to Work Plans and/or schedules.
- Activities undertaken in support of the Community Relations Plan.

Tasks completed during this period to fulfill each of these actions are summarized (by action) below.

Actions Taken During the Fourth Quarter to Achieve Compliance with the Consent Decree

The following actions were taken during September, October, and November of the Fourth Quarter of 2015:

- On September 8, 2015, AECOM submitted to the USEPA the Second Quarter GMZ Monitoring and System Performance Report.
- Phase 1 and Phase 2 air sparge/soil vapor extraction (AS/SVE) system air sampling of the SVE system effluent after being reactivated (switched to the pulse-on mode). SVE process air effluent sampling was conducted on September 22, 2015.
- Phase 1 and Phase 2 AS/SVE system air sampling of the SVE system effluent prior to the system being deactivated (switched to the pulse-off mode). SVE process air effluent sampling was conducted on November 20, 2015.

- Phase 2 SVE system water sample was collected from the Phase 2 Cell 4 and Cell 5 moisture separators. The moisture separator water samples were collected on November 20, 2015.

Summary of Sampling and Tests

- Three process air samples were collected from the Phase 1 AS/SVE system effluent during the September 22, 2015, and November 20, 2015, sampling events.
- Two process air samples were collected from the Phase 2 AS/SVE system effluent during the September 22, 2015 and November 20, 2015 sampling events.

Work Plans and Other Deliverables Completed In Accordance With the CD

- The Second Quarter GMZ Monitoring and System Performance Report, Area 9/10 Remedial Action (September 2015) was submitted in accordance with Section X, paragraph 30, of the CD and consistent with Section V of the Statement of Work (SOW).

Actions Scheduled for Next Quarter

The following actions are scheduled for the next quarter:

- Operation of the Phase 1 and Phase 2 AS/SVE systems will be in pulse-off mode (system not in operation) from November 20, 2015, to approximately January 18, 2016.
- Completion of the Fourth Quarter 2015 Groundwater Monitoring event in December 2015. The monitoring event will include the GMZ monitoring well network (which includes the Phase 1 AS/SVE performance monitoring network) and the Phase 2 AS/SVE performance monitoring network.
- Operation of the Phase 1 and Phase 2 AS/SVE systems will be in pulse-on mode (system in run mode) from approximately January 18, 2016, until approximately March 18, 2016.
- Completion of the First Quarter 2016 Groundwater Monitoring event in February 2016. The monitoring event will include the GMZ monitoring well network (which includes the Phase 1 AS/SVE performance monitoring network) and the Phase 2 AS/SVE performance monitoring network.

Percentage of Completion/Anticipated Delays

There are 39 specific deliverables or activities required to be completed as part of the CD. Some of these are ongoing activities and others, such as submittal of documents, require approval by the USEPA/Illinois Environmental Protection Agency to fulfill the requirements of the CD. To date, UTAS has completed their current obligations for 28 of the 39 items, or 72 percent of the CD requirements. There are currently no anticipated delays to the schedule.

Modifications to Work Plans/Schedules Proposed

None

Activities Undertaken In Support of Community Relations Plan

No activities are required with regard to the Community Relations Plan at this time.

